Final Fourth Quarter 2014 - Quarterly Groundwater Monitoring Report Inside Tunnel Wells

Red Hill Bulk Fuel Storage Facility
Joint Base Pearl Harbor-Hickam, Oahu, Hawaii

DOH Facility ID: 9-102271

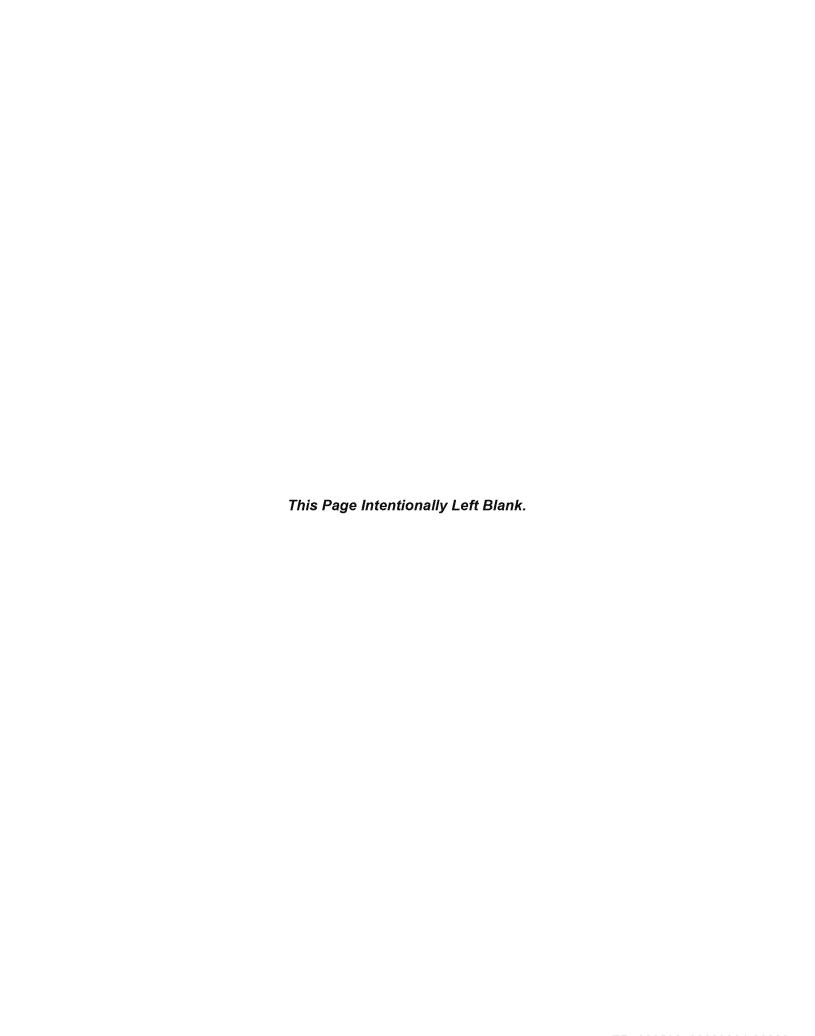
DOH Release ID: 990051, 010011, 020028, and 140010

January 2015

Department of the Navy Naval Facilities Engineering Command, Hawaii 400 Marshall Road JBPHH HI 96860-3139



Contract Number N62742-12-D-1853, CTO 0002



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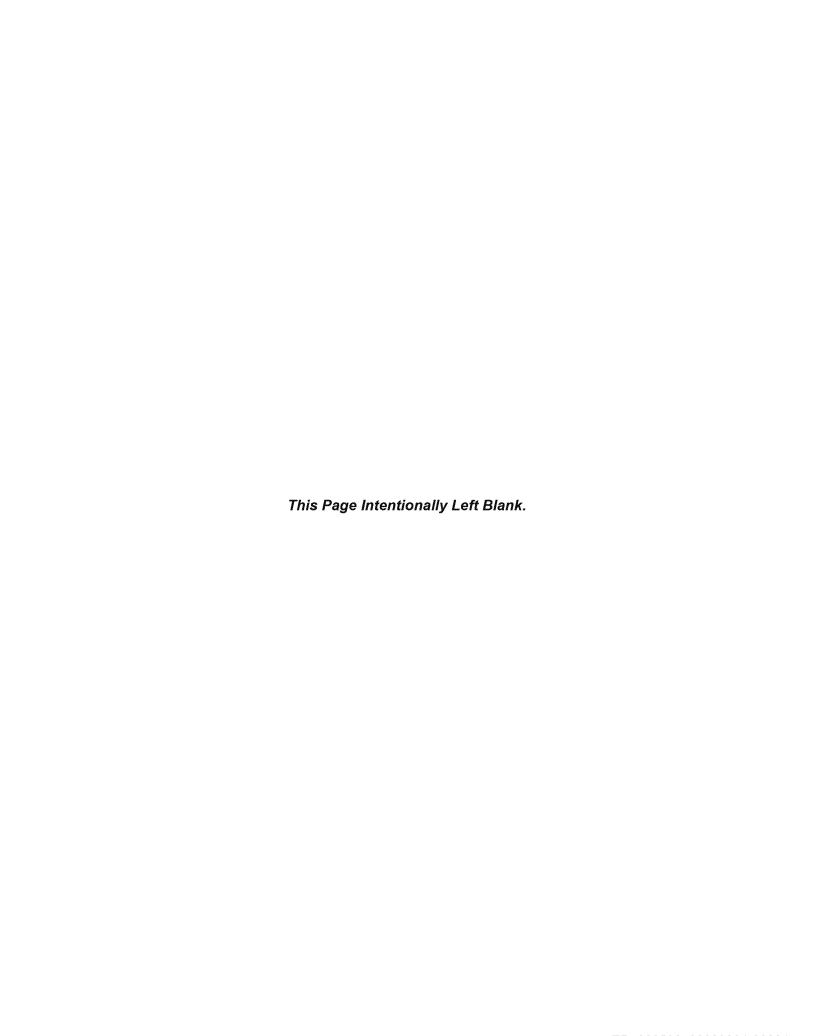
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Prepared under:

Contract Number N62742-12-D-1853, CTO 0002



FINAL FOURTH QUARTER 2014 - QUARTERLY GROUNDWATER MONITORING REPORT INSIDE TUNNEL WELLS RED HILL BULK FUEL STORAGE FACILITY

Long-Term Groundwater and Soil Vapor Monitoring Red Hill Bulk Fuel Storage Facility Joint Base Pearl Harbor-Hickam, Oahu, Hawaii

Prepared for:

Department of the Navy
Commanding Officer, Naval Facilities Engineering Command, Hawaii
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Prepared under:

Contract Number: N62742-12-D-1853 Contract Task Order: 0002

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ACRONYMS AND ABBREVIATIONS

ACRONYMS/

ABBREVIATIONS DEFINITION/MEANING

% percent

bgs below ground surface

COPC Contaminant of Potential Concern

DLNR State of Hawaii Department of Land and Natural Resources

DoD Department of Defense

DOH State of Hawaii Department of Health

DON Department of the Navy

EAL Environmental Action Level

EPA Environmental Protection Agency

ESI Environmental Science International, Inc.

F-76 Marine Diesel Fuel ID Identification

JBPHH Joint Base Pearl Harbor-Hickam

JP-5 Jet Fuel Propellant-5
JP-8 Jet Fuel Propellant-8
LCS Laboratory Control Sample

LCSD Laboratory Control Sample Duplicate

LOD Limit of Detection
LOQ Limit of Quantitation

µg/L micrograms per liter

MS Matrix Spike

MSD Matrix Spike Duplicate

NAVFAC Naval Facilities Engineering Command

NAVSUP FLC Naval Supply Systems Command Fleet Logistics Center

PAH Polycyclic Aromatic Hydrocarbons

PARCCS Precision, Accuracy, Representativeness, Completeness, Comparability,

and Sensitivity

pH hydrogen activity

RHSF Red Hill Bulk Fuel Storage Facility

RPD Relative Percent Difference
SAP Sampling and Analysis Plan
SSRBL Site-Specific Risk-Based Level
TEC The Environmental Company, Inc.
TPH-d Total Petroleum Hydrocarbons as diesel

TPH-g Total Petroleum Hydrocarbons as gasoline

U.S. United States of America
UST Underground Storage Tank
VOC Volatile Organic Compound

WP Work Plan

| Contract | Task | Order | 0002 |
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EXECUTIVE SUMMARY

This quarterly groundwater monitoring report presents the results of the fourth quarter 2014 groundwater sampling event, conducted on October 27 and 28, 2014, at the Red Hill Bulk Fuel Storage Facility [RHSF], Joint Base Pearl Harbor-Hickam [JBPHH], Hawaii. The RHSF is located in Halawa Heights on the Island of Oahu. There are 18 active and 2 inactive Underground Storage Tanks [USTs] located at the RHSF. The State of Hawaii Department of Health [DOH] Facility Identification [ID] number is 9-102271. The DOH Release ID numbers are 990051, 010011, 020028, and 140010.

The groundwater sampling was conducted as part of the long-term groundwater and soil vapor monitoring program at the RHSF and concurrent with release response activities initiated at Tank 5 in January 2014, under Naval Facilities Engineering Command [NAVFAC] Contract Number N62742-12-D-1853. The sampling was conducted in accordance with the approved 2012 Work Plan [WP]/Sampling and Analysis Plan [SAP] prepared by Environmental Science International, Inc. [ESI].

On October 27 and 28, 2014, ESI personnel collected groundwater samples from four monitoring wells at the RHSF (wells RHMW01, RHMW02, RHMW03, and RHMW05) and one sampling point at Red Hill Shaft (RHMW2254-01). A primary and duplicate sample were collected from well RHMW02. A summary of the analytical results is provided below.

- RHMW01 Total Petroleum Hydrocarbons as diesel fuel [TPH-d] (120 micrograms per liter [μg/L]) and lead (0.0976 μg/L) were the only analytes detected. The TPH-d concentration exceeded the DOH Environmental Action Level [EAL] for gross contamination, but was below the site-specific risk-based level [SSRBL] of 4,500 μg/L for Total Petroleum Hydrocarbons [TPH]. TPH-d concentrations in this well have shown an overall decreasing trend from a high of 1,500 μg/L in February 2005.
- RHMW02 TPH-d (2,000 μg/L in both primary and duplicate samples), Total Petroleum Hydrocarbons as gasoline [TPH-g] (57 and 53 μg/L), xylenes (0.32 and 0.29 μg/L), acenaphthene (0.53 μg/L in both primary and duplicate samples), ethylbenzene (0.15 and 0.14 μg/L), 1-methylnaphthalene (59 and 54 μg/L), 2-methylnaphthalene (43 and 36 μg/L), naphthalene (140 and 130 μg/L), and lead (0.165 μg/L in only the duplicate sample) were detected. TPH-d was detected at concentrations above the DOH EALs for both drinking water toxicity and gross contamination, but below half the SSRBL. The polycyclic aromatic hydrocarbons [PAHs] 1-methylnaphthalene, 2-methylnaphthalene, and naphthalene were detected at concentrations above the DOH EALs for both drinking water toxicity and gross contamination. The concentrations of 1-methylnaphthalene, 2-methylnaphthalene, and naphthalene increased to their highest levels in over 5 years. The concentrations of these PAHs are also significantly higher than they were in January 2014, when a high TPH-d spike was observed in this well.
- RHMW03 TPH-d (80 μg/L) was the only analyte detected. The concentration did not exceed the DOH EALs or the SSRBL.

- RHMW05 TPH-d (16 μ g/L) was the only analyte detected. The concentration did not exceed the DOH EALs or the SSRBL.
- RHMW2254-01 TPH-d (22 μg/L) and total lead (0.211 μg/L) were the only analytes detected. The concentrations did not exceed their respective DOH EAL.

During this quarterly event, the TPH-d concentration in RHMW01 exceeded the DOH EAL for gross contamination and the concentrations of TPH-d, 1-methylnaphthalene, 2-methylnaphthalene, and naphthalene in RHMW02 exceeded the DOH EALs for both gross contamination and drinking water toxicity. Groundwater contaminant concentrations in wells RHMW03, RHMW05, and RHMW2254-01 remained at low concentrations and did not change significantly from the previous event, or were not detected.

Concentrations of 1-methylnaphthalene, 2-methylnaphthalene, and naphthalene in RHMW02 have shown an increasing trend since March 5, 2014, but have remained below the historic maximums detected in the well. These PAH concentrations were an order of magnitude higher than they were during January 2014, when a TPH-d spike was observed but PAH concentrations remained below DOH EALs for gross contamination. Based on the groundwater monitoring data, TPH-d and PAH concentrations in RHMW02 appear to be decoupled. All other analytical results were generally consistent with historical data.

Based on the groundwater monitoring results and the reported release at Tank 5 in January 2014, continued groundwater monitoring at the wells inside the RHSF tunnel is recommended. The next quarterly event is tentatively scheduled for January 2015.

SECTION 1 – INTRODUCTION

This quarterly groundwater monitoring report presents the results of the fourth quarter 2014 groundwater sampling conducted on October 27 and 28, 2014, at the RHSF, JBPHH. The RHSF is located in Halawa Heights on the Island of Oahu. The purpose of the sampling is to (1) assess the condition of groundwater beneath the RHSF with respect to chemical constituents associated with jet fuel propellant and marine diesel fuel, and (2) to ensure the Navy remains in compliance with DOH UST release response requirements as described in Hawaii Administrative Rules Chapter 11-281 Subchapter 7, *Release Response Action* (DOH, 2013). The DOH Facility ID number for the RHSF is 9-102271. The DOH Release ID numbers are 990051, 010011, 020028, and 140010.

The groundwater sampling was conducted as part of the long-term groundwater and soil vapor monitoring program at the RHSF, under NAVFAC Contract Number N62742-12-D-1853. The sampling was conducted in accordance with the approved WP/SAP prepared by ESI (ESI, 2012).

1.1 SITE DESCRIPTION

The RHSF is located on federal government land (zoned F1- Military and Federal), located in Halawa Heights, approximately 2.5 miles northeast of Pearl Harbor (Figure 1). It is located on a low ridge on the western edge of the Koolau Mountain Range that divides Halawa Valley from Moanalua Valley. The RHSF is bordered on the north by Halawa Correctional Facility and private businesses, on the southwest by the United States of America [U.S.] Coast Guard reservation, on the south by residential neighborhoods, and on the east by Moanalua Valley. A quarry is located less than a quarter mile away to the northwest. The RHSF occupies 144 acres of land and the majority of the site is at an elevation of approximately 200 to 500 feet above mean sea level.

The RHSF contains 18 active and 2 inactive USTs that are operated by Naval Supply Systems Command Fleet Logistics Center [NAVSUP FLC] Pearl Harbor (formerly Fleet and Industrial Supply Center). Each UST has a capacity of approximately 12.5 million gallons. The RHSF is located approximately 100 feet above the basal aquifer. The USTs contain Jet Fuel Propellant-5 [JP-5], Jet Fuel Propellant-8 [JP-8], and Marine Diesel Fuel [F-76]. The current status of the USTs are summarized in Table 1.1.

Four groundwater monitoring wells (wells RHMW01, RHMW02, RHMW03, and RHMW05) and one sampling point at Red Hill Shaft (RHMW2254-01) are located within the RHSF lower access tunnel (Figure 2). Three groundwater monitoring wells (wells HDMW2253-03, OWDFMW01, and RHMW04) are located outside of the RHSF tunnel system. Monitoring data for the three wells located outside the tunnel are included in a separate report.

As noted, monitoring wells RHMW01, RHMW02, RHMW03, and RHMW05 are located inside the underground tunnels. Sampling point RHMW2254-01 is located inside the infiltration gallery

of the Department of the Navy [DON] drinking water supply Well 2254-01. The DON Well 2254-01 is located approximately 2,400 feet downgradient of the USTs and provides potable water to the JBPHH Water System, which serves approximately 65,200 military customers. NAVFAC Public Works Department operates the infiltration gallery and DON Well 2254-01.

TABLE 1.1
Current Status of the USTs
Red Hill Bulk Fuel Storage Facility
October 2014 Quarterly Monitoring Report

| | E 17 | | A |
|---------------------|-----------|----------|----------------------|
| Tank Identification | Fuel Type | Status | Capacity |
| F-1 | None | Inactive | 12.5 million gallons |
| F-2 | JP-8 | Active | 12.5 million gallons |
| F-3 | JP-8 | Active | 12.5 million gallons |
| F-4 | JP-8 | Active | 12.5 million gallons |
| F-5 | JP-8 | Active | 12.5 million gallons |
| F-6 | JP-8 | Active | 12.5 million gallons |
| F-7 | JP-5 | Active | 12.5 million gallons |
| F-8 | JP-5 | Active | 12.5 million gallons |
| F-9 | JP-5 | Active | 12.5 million gallons |
| F-10 | JP-5 | Active | 12.5 million gallons |
| F-11 | JP-5 | Active | 12.5 million gallons |
| F-12 | JP-5 | Active | 12.5 million gallons |
| F-13 | F-76 | Active | 12.5 million gallons |
| F-14 | F-76 | Active | 12.5 million gallons |
| F-15 | F-76 | Active | 12.5 million gallons |
| F-16 | F-76 | Active | 12.5 million gallons |
| F-17 | JP-5 | Active | 12.5 million gallons |
| F-18 | JP-5 | Active | 12.5 million gallons |
| F-19 | None | Inactive | 12.5 million gallons |
| F-20 | JP-5 | Active | 12.5 million gallons |

F-76 Marine Diesel Fuel

JP-5 Jet Fuel Propellant-5

JP-8 Jet Fuel Propellant-8

1.2 PHYSICAL SETTING

Climatological conditions in the area of the RHSF consist of warm to moderate temperatures and low to moderate rainfall. The RHSF is leeward of the prevailing northeasterly trade winds. The average annual precipitation is approximately 40 inches, which occurs mainly between November and April (State of Hawaii Department of Land and Natural Resources [DLNR], 1986). Annual pan evaporation is approximately 75 inches (DLNR, 1985). Average temperatures range from the low 60's to high 80's (degrees Fahrenheit) (Atlas of Hawaii, 1983).

Oahu consists of the eroded remnants of two shield volcanoes, Waianae and Koolau. The RHSF is located on the southwest flank of the Koolau volcanic shield. Lavas erupted during the shield-building phase of the volcano belong to the *Koolau Volcanic Series* (Stearns and Vaksvik, 1935). Following formation of the Koolau shield, a long period of volcanic quiescence occurred, during which the shield was deeply eroded. Following this erosional period, eruptive activity resumed. Lavas and pyroclastic material erupted during this period belong to the *Honolulu*

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Volcanic Series (Stearns and Vaksvik, 1935).

In the immediate area of the RHSF, Koolau Volcanic Series lavas dominate, although there are consolidated and unconsolidated non-calcareous deposits in the vicinity that consist of alluvium generated during erosion of the Koolau volcanic shield. South-southwest of the RHSF, and in isolated exposures to the west, are pyroclastic deposits formed during eruptions from three Honolulu Volcanic Series vents, Salt Lake, Aliamanu, and Makalapa (Stearns and Vaksvik, 1935). Based on established geology and records of wells drilled at the RHSF (Stearns and Vaksvik, 1938), the RHSF is underlain by Koolau Volcanic Series basalts. The area of the RHSF is classified as *Rock Land*, where 25-90% of the land surface is covered by exposed rock and there are only shallow soils (Foote, et al., 1972).

Groundwater in Hawaii exists in two principal aquifer types. The first and most important type, in terms of drinking water resources, is the basal aquifer. The basal aquifer exists as a lens of fresh water floating on and displacing seawater within the pore spaces, fractures, and voids of the basalt that forms the underlying mass of each Hawaiian island. In parts of Oahu, groundwater in the basal aquifer is confined by the overlying caprock and is under pressure. Waters that flow freely to the surface from wells that tap the basal aquifer are referred to as artesian.

The second type of aquifer is the caprock aquifer, which consists of various kinds of unconfined and semi-confined groundwater. Commonly, the caprock consists of a thick sequence of nearly impermeable clays, coral, and basalt, that separates the caprock aquifer from the basal aquifer. The impermeable nature of these materials and the artesian nature of the basal aquifer severely restrict the downward migration of groundwater from the upper caprock aquifer. However, in the area of the RHSF, there is no discernible caprock.

Groundwater in the area of the RHSF is part of the *Waimalu Aquifer System* of the *Pearl Harbor Aquifer Sector*. The aquifer is classified as a basal, unconfined, flank-type; and is currently used as a drinking water source. The aquifer is considered fresh, with less than 250 milligrams per liter of chloride, and is considered an irreplaceable resource with a high vulnerability to contamination (Mink and Lau, 1990).

The nearest drinking water supply well is DON Well 2254-01, located in the infiltration gallery within the RHSF lower tunnel. The DON Well 2254-01 is located approximately 2,400 feet hydraulically and topographically downgradient of the USTs (Figure 2).

1.3 BACKGROUND

The RHSF was constructed by the U.S. Government in the early 1940s. Twenty USTs and a series of tunnels were constructed. The USTs were constructed of steel, and in the past have stored DON special fuel oil, DON distillate, aviation gasoline, and motor gasoline (Environet, 2010). The tanks currently contain JP-5, JP-8, and F-76. The fueling system is a self-contained underground unit that was installed into native rock comprised primarily of basalt with some

interbedded tuffs and breccias (Environet, 2010). Each UST measures approximately 250 feet in height and 100 feet in diameter. The upper domes of the tanks lie at a depth varying between 100 feet and 200 feet below ground surface [bgs].

In 1998, Earth Tech conducted a Phase II Remedial Investigation/Feasibility Study for the Oily Waste Disposal Facility located within the RHSF. The study included the installation of well OWDFMW01 (which was originally MW08) (Earth Tech, 1999).

In February 2001, the DON installed groundwater monitoring well RHMW01 to monitor for contamination in the basal aquifer beneath the RHSF. Well RHMW01 was installed approximately 100 feet below grade within the lower access tunnel. The depth to water was measured at 86 feet below the tunnel floor at the time of the well completion. In February 2001, a groundwater sample was collected from the well. TPH and total lead were detected in the sample. Total lead was detected at a concentration above the DOH Tier 1 groundwater action level of 5.6 µg/L (The Environmental Company, Inc. [TEC], 2009; DOH, 2000).

In 2005, the RHSF groundwater monitoring program was initiated. It involved routine groundwater sampling of well RHMW01 and sampling point RHMW2254-01. Samples were collected in February, June, September, and December of 2005. Lead was detected at concentrations above the DOH Tier 1 action level of 5.6 µg/L in samples collected in February and June. The samples collected in February and June were not filtered prior to analysis, whereas the samples collected in September and December were filtered prior to analysis. Since the samples collected in February and June were not filtered prior to analysis, the lead results were not considered appropriate for a risk assessment (TEC, 2008).

Between June and September 2005, TEC installed three additional groundwater monitoring wells (wells RHMW02, RHMW03, and RHMW04) (TEC, 2008). Well RHMW04 was installed hydraulically upgradient of the USTs to provide background geochemistry information for water moving through the basal aquifer beneath the RHSF. Wells RHMW02 and RHMW03 were installed approximately 125 feet below grade within the RHSF lower tunnel and well RHMW04 was installed to a depth of approximately 300 feet bgs outside of the RHSF tunnels. In September 2005, groundwater samples were collected from the three newly installed groundwater monitoring wells (wells RHMW02, RHMW03, and RHMW04) along with the existing well RHMW01 and sampling point RHMW2254-01. The contaminants of potential concern [COPCs] with concentrations exceeding current DOH EALs are summarized below.

- RHMW01 TPH-d was detected at concentrations above the DOH EALs.
- RHMW02 TPH-g, TPH-d, naphthalene, trichloroethylene, 1-methylnaphthalene, and 2-methylnaphthalene were detected at concentrations above their respective DOH EAL.
- RHMW03 TPH-d was detected at concentrations above the DOH EALs.

In 2006, TEC installed dedicated sampling pumps in the four wells (wells RHWM01, RHWM02, RHMW03, and RHWM04) and one sampling point (RHMW2254-01). In July and December of

2006, groundwater samples were collected from the four wells and the sampling point. The COPCs with concentrations exceeding current DOH EALs are summarized below.

- RHMW01 TPH-d and naphthalene were detected at concentrations above their respective DOH EAL.
- RHMW02 TPH-g, TPH-d, and naphthalene were detected at concentrations above their respective DOH EAL.
- RHMW03 TPH-d was detected at concentrations above the DOH EALs.

In 2007, SSRBLs were established for TPH (4,500 μ g/L) and benzene (750 μ g/L) based on the solubility in water of JP-5 and JP-8 (TEC, 2007). Groundwater samples were collected from wells RHWM01, RHWM02, and RHMW03, and sampling point RHMW2254-01. Samples were collected in March, June, and September of 2007. The COPCs with concentrations exceeding current DOH EALs are summarized below.

- RHMW01 TPH-d was detected at concentrations above the DOH EALs, but below the SSRBL.
- RHMW02 TPH-g, TPH-d, naphthalene, 1-methylnaphthalene, and 2-methylnaphthalene were detected at concentrations above their respective DOH EAL. The TPH-d concentrations were below the SSRBL.
- RHMW03 TPH-d was detected at concentrations above the DOH EALs, but below the SSRBL.

In 2008, groundwater samples were collected from wells RHWM01, RHWM02, and RHMW03, and sampling point RHMW2254-01. Samples were collected in January, April, July, and October of 2008. The COPCs with concentrations exceeding current DOH EALs are summarized below. In addition, a groundwater protection plan (TEC, 2008) was prepared.

- RHMW01 TPH-d was detected at concentrations above the DOH EALs, but below the SSRBL.
- RHMW02 TPH-d, naphthalene, 1-methylnaphthalene, and 2-methylnaphthalene were
 detected at concentrations above their respective DOH EAL. The TPH-d concentrations
 detected in October 2008 were also above the SSRBL.
- RHMW03 TPH-d was detected at concentrations above the DOH EALs, but below the SSRBL.

In April 2009, groundwater monitoring well RHMW05 was installed downgradient of the USTs, within the lower access tunnel between RHMW01 and RHMW2254-01. It was installed to identify the extent of contamination hydraulically downgradient of the USTs. Well RHMW05 was added to the quarterly groundwater sampling program. In 2009, quarterly groundwater samples were collected from wells RHWM01, RHWM02, RHMW03, and RHMW05, and sampling point RHMW2254-01. Samples were collected in February, May, July, and October of 2009. The

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COPCs with concentrations exceeding current DOH EALs are summarized below. In addition, the Groundwater Protection Plan was revised to include well RHMW05.

- RHMW01 TPH-d and 1-methylnaphthalene were detected at concentrations above their respective DOH EAL. The TPH-d concentrations were below the SSRBL.
- RHMW02 TPH-d, naphthalene, 1-methylnaphthalene, and 2-methylnaphthalene were
 detected at concentrations above their respective DOH EAL. The TPH-d concentrations
 were below the SSRBL.
- RHMW03 TPH-d was detected at a concentration above the DOH EALs, but below the SSRBL.
- RHMW05 TPH-d was detected at a concentration above the DOH EALs, but below the SSRBI

In 2010, groundwater samples were collected from wells RHWM01, RHWM02, RHMW03, and RHMW05, and sampling point RHMW2254-01. Samples were collected in January, April, July, and October. The COPCs with concentrations exceeding current DOH EALs are summarized below.

- RHMW01 TPH-d was detected at concentrations above the DOH EALs, but below the SSRBL.
- RHMW02 TPH-g, TPH-d, naphthalene, and 1-methylnaphthalene were detected at concentrations above their respective DOH EAL. The TPH-d concentrations were below the SSRBL.
- RHMW03 TPH-d was detected at a concentration above the DOH EALs, but below the SSRBL.
- RHMW05 TPH-d was detected at a concentration above the DOH EALs, but below the SSRBL.

In 2011, quarterly groundwater samples were collected from wells RHWM01, RHWM02, RHMW03, and RHMW05, and sampling point RHMW2254-01. Samples were collected in January, April, July, and October. In a Fall 2011 update, the DOH EALs were revised. The drinking water toxicity EAL for TPH-d decreased from 210 to 190 μ g/L (DOH, 2011). The COPCs with concentrations exceeding current DOH EALs are summarized below.

- RHMW01 TPH-d was detected at concentrations above the DOH EALs, but below the SSRBL.
- RHMW02 TPH-d, naphthalene, indeno[1,2,3-cd]pyrene, and 1-methylnaphthalene were
 detected at concentrations above their respective DOH EAL. The TPH-d concentrations
 were below the SSRBL.

In 2012, quarterly groundwater samples were collected from wells RHWM01, RHWM02, RHMW03, and RHMW05, and sampling point RHMW2254-01. Samples were collected in

1-6

February, April, July, and November. The COPCs with concentrations exceeding current DOH EALs are summarized below.

- RHMW01 TPH-d was detected at concentrations above the DOH EALs, but below the SSRBL.
- RHMW02 TPH-d, TPH-g, naphthalene, 1-methylnaphthalene, and 2-methylnaphthalene were detected at concentrations above their respective DOH EAL. The TPH-d concentrations were below the SSRBL.

In 2013, quarterly groundwater samples were collected from wells RHWM01, RHWM02, RHMW03, and RHMW05, and sampling point RHMW2254-01. Samples were collected in January, April, July, and October. The COPCs with concentrations exceeding current DOH EALs are summarized below.

- RHMW01 TPH-d was detected at concentrations above the DOH EALs, but below the SSRBL.
- RHMW02 TPH-d, TPH-g, naphthalene, 1-methylnaphthalene, and 2-methylnaphthalene were detected at concentrations above their respective DOH EAL. The TPH-d concentrations were below the SSRBL.

In January 2014, quarterly groundwater samples were collected from wells RHMW01, RHMW02, RHMW03, RHMW05, and sampling point RHMW2254-01. The COPCs with concentrations exceeding current DOH EALs are summarized below.

- RHMW01 TPH-d was detected at a concentration above the DOH EALs, but below the SSRBL.
- RHMW02 TPH-d was detected at concentrations above both DOH EALs, and naphthalene
 and 1-methylnaphthalene were detected at concentrations above only the DOH EAL for
 drinking water toxicity. The TPH-d concentrations exceeded the SSRBL.

Between January and April 2014, additional groundwater sampling was conducted at wells RHMW01, RHMW02, RHMW05, and sampling point RHMW2254-01 in response to a reported release from Tank 5. The COPCs with concentrations exceeding current DOH EALs are summarized below.

- RHMW01 TPH-d was detected at concentrations above the DOH EALs, but below the SSRBL.
- RHMW02 TPH-d was detected at concentrations above both DOH EALs, and naphthalene
 and 1-methylnaphthalene were detected at concentrations above only the DOH EAL for
 drinking water toxicity. The TPH-d concentrations were below the SSRBL.

In April 2014, quarterly groundwater samples were collected from wells RHMW01, RHMW02, RHMW03, RHMW05, and sampling point RHMW2254-01. The COPCs with concentrations exceeding current DOH EALs are summarized below.

 RHMW02 – TPH-d and naphthalene were detected at concentrations above both DOH EALs. 1-Methylnaphthalene was detected at concentrations above only the DOH EAL for drinking water toxicity. The TPH-d concentrations were below the SSRBL.

In May and June 2014, additional groundwater sampling was conducted at wells RHMW01, RHMW02, RHMW05, and sampling point RHMW2254-01 in response to a reported release from Tank 5. The COPCs with concentrations exceeding current DOH EALs are summarized below.

 RHMW02 – TPH-d, naphthalene, and 1-methylnaphthalene were detected at concentrations above their respective DOH EAL. The TPH-d concentrations were below the SSRBL.

In July 2014, quarterly groundwater samples were collected from wells RHMW01, RHMW02, RHMW03, RHMW05, and sampling point RHMW2254-01. The COPCs with concentrations exceeding current DOH EALs are summarized below.

 RHMW02 – TPH-d, 1-methylnaphthalene, and naphthalene were detected at concentrations above their respective DOH EAL. 2-Methylnaphthalene was detected at concentrations above only the DOH EAL for gross contamination. The TPH-d concentrations were below the SSRBL.

1.3.1 Previous Reports

The following groundwater monitoring reports were previously submitted to the DOH:

- 1. Groundwater Sampling Report, First Quarter 2005 (submitted April 2005).
- 2. Groundwater Sampling Report, Second Quarter 2005 (submitted August 2005).
- 3. Groundwater Sampling Report, Third Quarter 2005 (submitted November 2005).
- 4. Groundwater Sampling Report, Fourth Quarter 2005 (submitted February 2006).
- 5. Groundwater Monitoring Results, July 2006 (submitted September 2006).
- Groundwater Monitoring Results, December 2006 (submitted January 2007).
- 7. Groundwater Monitoring Results, March 2007 (submitted May 2007).
- 8. Groundwater Monitoring Results, June 2007 (submitted August 2007).
- Groundwater Monitoring Results, September 2007 (submitted October 2007).
- 10. Groundwater Monitoring Report, January 2008 (submitted March 2008).
- 11. Groundwater Monitoring Report, April 2008 (submitted May 2008).

- 12. Groundwater Monitoring Report, July 2008 (submitted October 2008).
- 13. Groundwater Monitoring Report, October and December 2008 (submitted February 2009).
- 14. Groundwater Monitoring Report, February 2009 (submitted May 2009).
- 15. Groundwater Monitoring Report, May 2009 (submitted July 2009).
- 16. Groundwater Monitoring Report, July 2009 (submitted September 2009).
- 17. Groundwater Monitoring Report, October 2009 (submitted December 2009).
- 18. Groundwater Monitoring Report, January, February, and March 2010 (submitted April 2010).
- 19. Groundwater Monitoring Report, April 2010 (submitted May 2010).
- 20. Groundwater Monitoring Report, July 2010 (submitted August 2010).
- 21. Groundwater Monitoring Report, October 2010 (submitted December 2010).
- 22. Groundwater Monitoring Report, January 2011 (submitted March 2011).
- 23. Groundwater Monitoring Report, April 2011 (submitted June 2011).
- 24. Groundwater Monitoring Report, July 2011 (submitted September 2011).
- 25. Groundwater Monitoring Report, October 2011 (submitted December 2011).
- 26. Groundwater Monitoring Report, January-February 2012 (submitted March 2012).
- 27. Groundwater Monitoring Report, April 2012 (Submitted July 2012).
- 28. Groundwater Monitoring Report, October 2012 (Submitted January 2013).
- 29. Groundwater Monitoring Report, January 2013 (Submitted April 2013).
- 30. Groundwater Monitoring Report, April 2013 (Submitted July 2013).
- 31. Groundwater Monitoring Report, July 2013 (Submitted September 2013).
- 32. Groundwater Monitoring Report, October 2013 (Submitted January 2014).
- 33. Groundwater Sampling Report for Additional Sampling, January 2014 (submitted January 2014).
- 34. Groundwater Monitoring Report, January 2014 (Submitted April 2014).
- 35. Groundwater Sampling Report for Tank 5 Release Response on March 5 and 6, 2014 (submitted March 2014).
- 36. Groundwater Sampling Report for Tank 5 Release Response on March 10, 2014 (submitted March 2014).
- 37. Groundwater Sampling Report for Tank 5 Release Response on March 25 and 26, 2014 (submitted April 2014).

- 38. Groundwater Sampling Report for Tank 5 Release Response on April 7, 2014 (submitted April 2014).
- 39. Groundwater Monitoring Report, April 2014 (Submitted June 2014).
- 40. Groundwater Sampling Report for Tank 5 Release Response on May 27 and 28, 2014 (submitted June 2014).
- 41. Groundwater Sampling Report for Tank 5 Release Response on June 23 and 24, 2014 (submitted July 2014).
- 42. Groundwater Monitoring Report, July 2014 (Submitted September 2014).

SECTION 2 – GROUNDWATER SAMPLING

On October 27 and 28, 2014, ESI personnel collected groundwater samples from four monitoring wells at the RHSF (wells RHMW01, RHMW02, RHMW03, and RHMW05) and one sampling point at Red Hill Shaft (RHMW2254-01). The samples were collected in accordance with the 2012 WP/SAP. The WP/SAP is consistent with DOH UST release response requirements (DOH, 2000); DON Procedure I-C-3, *Monitoring Well Sampling* (DON, 2007); and the RHSF Groundwater Protection Plan (TEC, 2008). Prior to purging and sampling, the depth to groundwater and the depth to the bottoms of the wells were measured using a Geotech oil/water interface probe. No measurable product, sheen, or petroleum hydrocarbon odor was detected in any of the wells.

2.1 GROUNDWATER SAMPLING

Prior to collecting groundwater samples, the monitoring wells were purged of standing water in the well casings. Each well contains a dedicated bladder pump which was used to purge the well and to collect samples. To operate the pump, a portable air compressor with an in-line filter was connected to a QED MP50 MicroPurge® Basics Controller box, which was then connected to the pump. The compressor was turned on to power the pump and the controller was used to adjust the pumping rate to less than one liter of water per minute.

Water quality parameters were monitored on a periodic basis during well purging. The water quality parameters that were measured included hydrogen activity [pH], temperature, conductivity, dissolved oxygen, and oxidation reduction potential. The water quality parameters were evaluated to assess whether the natural characteristics of the aquifer formation water were present within the monitoring wells before collecting the samples. At least four readings were collected during the purging process. Purging was considered complete when at least three consecutive water quality measurements stabilized within approximately 10%. The readings were recorded on groundwater monitoring logs. The groundwater monitoring logs are included in Appendix A. In addition, field notes were taken to document the sampling event. The field notes are included in Appendix B.

When the water quality parameters stabilized, groundwater samples were collected from the wells using the bladder pumps. The groundwater samples were collected no more than two hours after purging was completed to decrease groundwater interaction with the monitoring well casing and atmosphere. Prior to collecting the sample, the water level in the monitoring wells was measured and recorded to ensure that excessive drawn down had not occurred. The groundwater samples were collected at flow rates of approximately 0.17 to 0.5 liters per minute. Samples collected for dissolved lead analysis were filtered in the field using new, 0.45-micron filters.

All samples were labeled and logged on the Sample Inventory Log, placed in Ziploc™ bags and sealed, custody sealed, sealed with tape, placed in a cooler with wet ice, and logged onto the Chain-of-Custody form. The samples were labeled and logged in accordance with Project

Procedures Manual Standard Operating Procedure III-E, Record Keeping, Sample Labeling, and Chain-of-Custody Procedures (DON, 2007). All samples were shipped under Chain-of-Custody to the analytical laboratory and analyzed for the COPCs as described in Section 2.2.

2.2 ANALYTICAL RESULTS

The samples were analyzed for TPH-d using U.S. Environmental Protection Agency [EPA] Method 8015M, TPH-g and Volatile Organic Compounds [VOCs] using EPA Method 8260B, PAHs using EPA Method 8270C SIM, dissolved lead using EPA Method 6020, and total lead using EPA Method 200.8. The sample collected from sampling point RHMW2254-01 was analyzed for total lead (unfiltered) as DON Well 2254-01 is a drinking water supply well. A copy of the laboratory report is included as Appendix C.

Analytical results were compared to the DOH EALs for drinking water toxicity and gross contamination. Analytical results for wells RHMW01, RHMW02, RHMW03, and RHMW05 were also compared to the SSRBLs for TPH (4,500 μ g/L) and benzene (750 μ g/L), established in the RHSF Final Groundwater Protection Plan (TEC, 2008). The results of the fourth quarter groundwater sampling event are summarized in Table 2.1 and described below.

- RHMW01 TPH-d (120 µg/L) and lead (0.0976 µg/L) were the only analytes detected. The
 TPH-d concentration exceeded the DOH EAL for gross contamination, but was below the
 SSRBL of 4,500 µg/L for TPH. None of the other chemical constituents analyzed for were
 detected at concentrations above their respective DOH EAL.
- RHMW02 TPH-d (2,000 μg/L), TPH-g (57 and 53 μg/L), xylenes (0.32 and 0.29 μg/L), acenaphthene (0.53 μg/L), ethylbenzene (0.15 and 0.14 μg/L), 1-methylnaphthalene (59 and 54 μg/L), 2-methylnaphthalene (43 and 36 μg/L), and naphthalene (140 and 130 μg/L) were detected in both the primary and duplicate samples collected. Lead (0.165 μg/L) was only detected in the duplicate sample. TPH-d, 1-methylnaphthalene, 2-methylnaphthalene, and naphthalene were detected at concentrations above their respective DOH EALs for both drinking water toxicity and gross contamination. However, the TPH-d concentrations did not exceed the SSRBL.
- RHMW03 TPH-d (80 μg/L) was the only analyte detected. The concentration did not exceed the DOH EAL or the SSRBL.
- RHMW05 TPH-d (16 μg/L) was the only analyte detected. The concentration did not exceed the DOH EAL or the SSRBL.
- RHMW2254-01 TPH-d (22 μg/L) and total lead (0.211 μg/L) were the only analytes detected. The concentrations did not exceed their respective DOH EAL.

2.3 GROUNDWATER CONTAMINANT TRENDS

The historical groundwater contaminant concentrations for COPCs that exceeded the DOH EALs or SSRBLs are illustrated in Appendix D. A table of cumulative historical groundwater

results is included as Appendix E. A summary of groundwater contaminant trends is provided below.

- RHMW01 COPCs detected during this round of quarterly sampling are consistent with the historical data for RHMW01. TPH-d has historically been detected at concentrations above the DOH EAL for both drinking water toxicity and gross contamination. TPH-d concentrations have shown an overall decreasing trend from a high of 1,500 µg/L in February 2005. The TPH-d concentration exceeded the DOH EAL for gross contamination for the first time since April 7, 2014.
- RHMW02 COPCs detected during this round of quarterly sampling are consistent with the historical data for RHMW02. TPH-g, TPH-d, 1-methylnaphthalene, 2-methylnaphthalene, and naphthalene have historically been detected at concentrations above the DOH EALs. During the October 2014 event, concentrations of TPH-d, 1-methylnaphthalene, 2-methylnaphthalene, and naphthalene approximately doubled from the previous event in July 2014 and remained above DOH EALs. The concentrations of 1-methylnaphthalene, 2-methylnaphthalene, and naphthalene increased to their highest levels in over 5 years. The concentrations of these PAHs are also significantly higher than they were in January 2014, when a high TPH-d spike was observed in this well. The concentrations of TPH-g remained below the DOH EALs for gross contamination and drinking water toxicity and were comparable to the concentrations detected during the previous event. Trichloroethylene was detected once in RHMW02 in September 2005 in the primary sample at a concentration above the DOH EAL for drinking water toxicity; however, trichloroethylene was not detected in the duplicate sample, and this may have been an anomalous result.
- RHMW03 COPCs detected during this round of quarterly sampling are consistent with the
 historical data for RHMW03. TPH-d has historically been detected at concentrations above
 the DOH EALs; however, it has not been detected at concentrations above the DOH EALs
 since October 2010.
- RHMW05 COPCs detected during this round of quarterly sampling are consistent with the
 historical data for RHMW05. TPH-d has historically been detected in RHMW05 at
 concentrations above the DOH EAL for both drinking water toxicity and gross contamination;
 however, it has not been detected at concentrations above the DOH EALs since January
 2010.
- RHMW2254-01 COPCs detected during this round of quarterly sampling are consistent
 with the historical data for RHMW2254-01. Although the method reporting limits for TPH-d
 exceeded one or both DOH EALs for drinking water toxicity and gross contamination
 between May 2009 and July 2010, TPH-d was last detected in RHMW2254-01 at a
 concentration above the DOH EAL for gross contamination in January 2008.

2.4 WASTE DISPOSAL

The purged groundwater and decontamination water generated during sampling of the inside tunnel was placed in a 55-gallon drum along with the purged water and decontamination water from the outside tunnel wells. The drum is currently stored onsite at ADIT 3 on top of a

secondary containment spill pallet and covered by a tarp. There is a non-hazardous label affixed to the drum with all pertinent information relating to its generation. The drum will be used for future sampling events and will be properly disposed of once it has been filled.

Contract No. N62742-12-D-1853 Contract Task Order 0002

TABLE 2.1

Analytical Results for Groundwater Sampling (October 27 and 28, 2014) Red Hill Bulk Fuel Storage Facility October 2014 Quarterly Monitoring Report

| | | DOH E | ALs | | RHMW: | 2254-01 | ES117) | | | RHV | VM01 (ES1 | 13) | | | RH | MW02 (ES | 114) | | | RHV | VM03 (ES | 116) | | | RHN | /W05 (ES | 118) | |
|------------------------|---|-----------------|------------------|--------------|--------------|--------------|----------|----------------|--------------------|--------------|-----------|----------------|-------------|--------------------|--|--------------|----------------|-------------|--------------|------------|----------|-------|--------------|--------------|------------|--------------|----------|-------------|
| Method | Chemical | Drinking Water | Gross | Results | o | LOQ | LOD | DL | Results | 0 | LOQ | LOD | DL | Results | a | Log | LOD | DL | Results | o | LOQ | LOD | DL | Results | 0 | LOQ | LOD | DL |
| EDA 901ED | TDU d | Toxicity | Contamination | | | | | | | | | | | | | | | 11 | | HD,J | | | 11 | 16 | 1 | | | |
| EPA 8015B EPA 8260B | TPH-a | 190 100 | 100 | 22 N.D. | HD,J U | 25 50 | 12 30 | 11 26 | 120 N.D. | HD,J U | 25 50 | 12 30 | 11 26 | 2,000 57 | HD,J | 25 50 | 12 30 | 26 | 80 N.D. | U U | 25 50 | 30 | 26 | N.D. | HD,J U | 25 50 | 12 30 | 11 26 |
| | Acenaphthene | 370 | 20 | N.D. | Ü | 0.19 | 0.049 | 0.026 | N.D. | Ū | 0.21 | 0.052 | 0.028 | 0.53 | † | 0.19 | 0.047 | 0.025 | N.D. | Ü | 0.21 | 0.054 | 0.029 | N.D. | Ü | 0.19 | 0.048 | 0.026 |
| | Acenaphthylene | 240 | 2,000 | N.D. | U | 0.19 | 0.049 | 0.043 | N.D. | U | 0.21 | 0.052 | 0.046 | N.D. | U | 0.19 | 0.047 | 0.042 | N.D. | U | 0.21 | 0.054 | 0.048 | N.D. | U | 0.19 | 0.048 | 0.043 |
| | Anthracene | 1,800 | 22 | N.D. | U | 0.19 | 0.049 | 0.028 | N.D. | U | 0.21 | 0.052 | 0.030 | N.D. | U | 0.19 | 0.047 | 0.027 | N.D. | U | 0.21 | 0.054 | 0.031 | N.D. | U | 0.19 | 0.048 | 0.028 |
| | Benzo[a]anthracene | 0.092 1,500 | 4.7 0.13 | N.D. | U | 0.19 | 0.049 | 0.032 0.080 | N.D. N.D. | U | 0.21 | 0.052 0.10 | 0.034 | N.D. N.D. | U | 0.19 0.19 | 0.047 0.095 | 0.031 | N.D. | U | 0.21 | 0.054 | 0.035 | N.D. | U | 0.19 0.19 | 0.048 | 0.031 |
| | Benzo[g,h,i]perylene Benzo[a]pyrene | 0.2 | 0.13 | N.D. | u | 0.19 | 0.097 | 0.000 | N.D. | - | 0.21 | 0.10 | 0.003 | N.D. | U | 0.19 | 0.093 | 0.076 | N.D. | U | 0.21 | 0.054 | 0.000 | N.D. | U | 0.19 | 0.098 | 0.079 |
| | Benzo[b]fluoranthene | 0.092 | 0.75 | N.D. | Ű | 0.19 | 0.049 | 0.017 | N.D. | Ū | 0.21 | 0.052 | 0.018 | N.D. | Ü | 0.19 | 0.047 | 0.017 | N.D. | Ü | 0.21 | 0.054 | 0.019 | N.D. | Ü | 0.19 | 0.048 | 0.017 |
| | Benzo[k]fluoranthene | 0.92 | 0.4 | N.D. | U | 0.19 | 0.049 | 0.030 | N.D. | U | 0.21 | 0.052 | 0.032 | N.D. | U | 0.19 | 0.047 | 0.029 | N.D. | U | 0.21 | 0.054 | 0.033 | N.D. | U | 0.19 | 0.048 | 0.030 |
| EPA 8270C | Chrysene | 9.2 | 1 2.50 | N.D. | U | 0.19 | 0.049 | 0.024 | N.D. | U | 0.21 | 0.052 | 0.026 | N.D. | U | 0.19 | 0.047 | 0.024 | N.D. | U | 0.21 | 0.054 | 0.027 | N.D. | U | 0.19 | 0.048 | 0.024 |
| | Dibenzo[a,h]anthracene Fluoranthene | 0.0092 1,500 | 0.52 130 | N.D. | U | 0.19 | 0.049 | 0.046 0.045 | N.D. | U | 0.21 | 0.052 0.052 | 0.049 | N.D. N.D. | U | 0.19 0.19 | 0.047 0.047 | 0.045 | N.D. | U | 0.21 | 0.054 | 0.051 | N.D. | U | 0.19 0.19 | 0.048 | 0.046 |
| | Fluorene | 240 | 950 | N.D. | Ü | 0.19 | 0.049 | 0.043 | N.D. | Ū | 0.21 | 0.052 | 0.044 | N.D. | U | 0.19 | 0.047 | 0.040 | N.D. | Ü | 0.21 | 0.054 | 0.046 | N.D. | Ü | 0.19 | 0.048 | 0.043 |
| | Indeno[1,2,3-cd]pyrene | 0.092 | 0.095 | N.D. | Ū | 0.19 | 0.049 | 0.021 | N.D. | Ū | 0.21 | 0.052 | 0.022 | N.D. | Ü | 0.19 | 0.047 | 0.020 | N.D. | Ū | 0.21 | 0.054 | 0.023 | N.D. | Ü | 0.19 | 0.048 | 0.020 |
| | 1-Methylnaphthalene | 4.7 | 10 | N.D. | U | 0.19 | 0.097 | 0.050 | N.D. | U | 0.21 | 0.10 | 0.054 | 59 | | 3.8 | 1.9 | 0.98 | N.D. | U | 0.21 | 0.011 | 0.055 | N.D. | U | 0.19 | 0.096 | 0.050 |
| | 2-Methylnaphthalene | 24 | 10 | N.D. | U | 0.19 | 0.049 | 0.045 | N.D. | U | 0.21 | 0.052 | 0.048 | 43 | | 3.8 | 0.95 | 0.88 | N.D. | U | 0.21 | 0.054 | 0.050 | N.D. | U | 0.19 | 0.048 | 0.045 |
| | Naphthalene Phenanthrene | 17 240 | 21 410 | N.D. N.D. | U | 0.19 0.19 | 0.049 | 0.033 0.026 | N.D. N.D. | U | 0.21 | 0.052 0.052 | 0.035 | 140 N.D. | U | 3.8 0.19 | 0.95 0.047 | 0.64 | N.D. | U | 0.21 | 0.054 | 0.036 | N.D. N.D. | U | 0.19 0.19 | 0.048 | 0.033 |
| | Pyrene | 180 | 68 | N.D. | U | 0.19 | 0.049 | 0.020 | N.D. | U | 0.21 | 0.052 | 0.026 | N.D. | U | 0.19 | 0.047 | 0.026 | N.D. | U | 0.21 | 0.054 | 0.029 | N.D. | U | 0.19 | 0.048 | 0.020 |
| | 1,1,1,2-Tetrachloroethane | 0.52 | 50,000 | N.D. | Ü | 1 | 0.5 | 0.4 | N.D. | Ü | 1 | 0.5 | 0.4 | N.D. | Ü | 1 | 0.5 | 0.4 | N.D. | Ü | 1 | 0.5 | 0.4 | N.D. | Ü | 1 | 0.5 | 0.4 |
| | 1,1,2,2-Tetrachloroethane ¹ | 0.067 | 500 | N.D. | U | 1 | 0.5 | 0.41 | N.D. | U | 1 | 0.5 | 0.41 | N.D. | U | 1 | 0.5 | 0.41 | N.D. | U | 1 | 0.5 | 0.41 | N.D. | U | 1 | 0.5 | 0.41 |
| | 1,1,1-Trichloroethane | 200 | 970 | N.D. | U | 5 | 0.5 | 0.3 | N.D. | U | 5 | 0.5 | 0.3 | N.D. | U | 5 | 0.5 | 0.3 | N.D. | U | 5 | 0.5 | 0.3 | N.D. | U | 5 | 0.5 | 0.3 |
| | 1,1,2-Trichloroethane 1,1-Dichloroethane | 5 2.4 | 50,000 50,000 | N.D. | U | 5 | 0.5 | 0.38 0.28 | N.D. N.D. | U | 5 | 0.5 0.5 | 0.38 | N.D. N.D. | U | 1 5 | 0.5 0.5 | 0.38 | N.D. | U | 5 | 0.5 | 0.38 0.28 | N.D. N.D. | U | 1 5 | 0.5 | 0.38 |
| | 1,1-Dichloroethylene | 7 | 1,500 | N.D. | Ü | 1 | 0.5 | 0.28 | N.D. | U | 1 | 0.5 | 0.28 | N.D. | U | 1 | 0.5 | 0.43 | N.D. | U | 1 | 0.5 | 0.28 | N.D. | U | 1 | 0.5 | 0.43 |
| | 1,2,3-Trichloropropane ¹ | 0.6 | 50,000 | N.D. | Ü | 5 | 1 | 0.64 | N.D. | Ū | 5 | 1 | 0.64 | N.D. | Ü | 5 | 1 | 0.64 | N.D. | Ü | 5 | 1 | 0.64 | N.D. | Ü | 5 | 1 | 0.64 |
| | 1,2,4-Trichlorobenzene | 70 | 3,000 | N.D. | U | 5 | 1 | 0.5 | N.D. | U | 5 | 1 | 0.5 | N.D. | U | 5 | 1 | 0.5 | N.D. | U | 5 | 1 | 0.5 | N.D. | U | 5 | 1 | 0.5 |
| | 1,2-Dibromo-3- chloropropane | 0.04 | 10 | N.D. | U | 10 | 2 | 1.2 | N.D. | U | 10 | 2 | 1.2 | N.D. | U | 10 | 2 | 1.2 | N.D. | U | 10 | 2 | 1.2 | N.D. | U | 10 | 2 | 1.2 |
| | 1,2-Dibromoethane ¹ 1,2-Dichlorobenzene | 0.04 600 | 50,000 10 | N.D. | U | 1 1 | 0.5 | 0.24 0.46 | N.D. N.D. | U | 1 | 0.5 0.5 | 0.24 | N.D. | U | 1 1 | 0.5 | 0.24 | N.D. | U | 1 | 0.5 | 0.24 | N.D. | U | 1 | 0.5 | 0.24 |
| | 1,2-Dichloroethane | 0.15 | 7,000 | N.D. | Ü | 1 | 0.5 | 0.24 | N.D. | U,IH | 1 | 0.5 | 0.40 | N.D. | U,IH | 1 | 0.5 | 0.24 | N.D. | U,IH | 1 | 0.5 | 0.40 | N.D. | U | 1 | 0.5 | 0.24 |
| | 1,2-Dichloropropane | 5 | 10 | N.D. | Ū | 5 | 0.5 | 0.42 | N.D. | U | 5 | 0.5 | 0.42 | N.D. | U | 5 | 0.5 | 0.42 | N.D. | U | 5 | 0.5 | 0.42 | N.D. | Ü | 5 | 0.5 | 0.42 |
| | 1,3-Dichlorobenzene | 180 | 5 | N.D. | U | 1 | 0.5 | 0.4 | N.D. | U | 1 | 0.5 | 0.4 | N.D. | U | 1 | 0.5 | 0.4 | N.D. | U | 1 | 0.5 | 0.4 | N.D. | U | 1 | 0.5 | 0.4 |
| | 1,3-Dichloropropene (total of cis/trans) 1 | 0.43 | 50,000 | N.D. | U | 1 | 0.5 | 0.25 | N.D. | U | 1 | 0.5 | 0.25 | N.D. | U | 1 | 0.5 | 0.25 | N.D. | υ | 1 | 0.5 | 0.25 | N.D. | U | 1 | 0.5 | 0.25 |
| 1 | 1,4-Dichlorobenzene | 75 | 5 | N.D. | U | 1 | 0.5 | 0.43 | N.D. | U | 1 | 0.5 | 0.43 | N.D. | U | 1 | 0.5 | 0.43 | N.D. | U | 1 | 0.5 | 0.43 | N.D. | U | 1 | 0.5 | 0.43 |
| | Acetone | 22,000 | 20,000 | N.D. | U,ICH | 20 | 10 | 6 | N.D. | U,ICH | 20 | 10 | 6 | N.D. | U,ICH | 20 | 10 | 6 | N.D. | U,ICH | 20 | 10 | 6 | N.D. | U,ICH | 20 | 10 | 6 |
| | Benzene | 5 | 170 | N.D. | U | 1 | 0.5 | 0.14 | N.D. | U | 1 | 0.5 | 0.14 | N.D. | U | 1 | 0.5 | 0.14 | N.D. | U | 1 | 0.5 | 0.14 | N.D. | U | 1 | 0.5 | 0.14 |
| | Bromodichloromethane ¹ | 0.12 | 50,000 | N.D. | U | 5 | 0.5 | 0.21 | N.D. | U | 5 | 0.5 | 0.21 | N.D. | U | 5 | 0.5 | 0.21 | N.D. | U | 5 | 0.5 | 0.21 | N.D. | U | 5 | 0.5 | 0.21 |
| | Bromoform Bromomethane | 80 8.7 | 510 50,000 | N.D. | U | 10 20 | 5 | 0.5 3.9 | N.D. N.D. | U | 10 20 | 1 5 | 0.5 3.9 | N.D. | U | 10 20 | 5 | 0.5 3.9 | N.D. | U | 10 20 | 1 | 0.5 3.9 | N.D. | U | 10 20 | 1 5 | 0.5 3.9 |
| EPA 8260B | Carbon Tetrachloride | 5 | 520 | N.D. | Ü | 1 | 0.5 | 0.23 | N.D. | U | 1 | 0.5 | 0.23 | N.D. | U | 1 | 0.5 | 0.23 | N.D. | U | 1 | 0.5 | 0.23 | N.D. | U | 1 | 0.5 | 0.23 |
| | Chlorobenzene | 100 | 50 | N.D. | Ü | 5 | 0.5 | 0.17 | N.D. | Ū | 5 | 0.5 | 0.17 | N.D. | Ü | 5 | 0.5 | 0.17 | N.D. | Ű | 5 | 0.5 | 0.17 | N.D. | Ü | 5 | 0.5 | 0.17 |
| | Chloroethane | 21,000 | 16 | N.D. | U | 10 | 5 | 2.3 | N.D. | U | 10 | 5 | 2.3 | N.D. | U | 10 | 5 | 2.3 | N.D. | U | 10 | 5 | 2.3 | N.D. | U | 10 | 5 | 2.3 |
| | Chloroform | 70 | 2,400 | N.D. | U | 5 | 0.5 | 0.46 | N.D. | U | 5 | 0.5 | 0.46 | N.D. | U | 5 | 0.5 | 0.46 | N.D. | U | 5 | 0.5 | 0.46 | N.D. | U | 5 | 0.5 | 0.46 |
| | Chloromethane ¹ cis-1,2-Dichloroethylene | 1.8 70 | 50,000 50,000 | N.D. | U, IJ U | 10 | 0.5 | 1.8 0.48 | N.D. N.D. | U, IJ U | 10 | 2 0.5 | 1.8 0.48 | N.D. | U, IJ | 10 | 0.5 | 1.8 0.48 | N.D. | U, IJ U | 10 | 0.5 | 1.8 0.48 | N.D. | U, IJ U | 10 | 0.5 | 1.8 0.48 |
| | Dibromochloromethane ¹ | 0,16 | 50,000 | N.D. | Ü | 1 | 0.5 | 0.46 | N.D. | U | 1 1 | 0.5 | 0.46 | N.D. | U | 1 | 0.5 | 0.46 | N.D. | U | 1 | 0.5 | 0.46 | N.D. | U | 1 | 0.5 | 0.46 |
| | Ethylbenzene | 700 | 30 | N.D. | Ü | 1 | 0.5 | 0.14 | N.D. | Ū | 1 | 0.5 | 0.14 | 0.15 | J | 1 | 0.5 | 0.14 | N.D. | Ü | 1 | 0.5 | 0.14 | N.D. | Ü | 1 | 0.5 | 0.14 |
| | Hexachlorobutadiene | 0.86 | 6 | N.D. | U | 1 | 0.5 | 0.32 | N.D. | U | 1 | 0.5 | 0.32 | N.D. | U | 1 | 0.5 | 0.32 | N.D. | U | 1 | 0.5 | 0.32 | N.D. | U | 1 | 0.5 | 0.32 |
| | Methyl ethyl ketone (2- | 7,100 | 8,400 | N.D. | U | 10 | 5.0 | 2.2 | N.D. | U,ICH | 10 | 5.0 | 2.2 | N.D. | U,ICH | 10 | 5.0 | 2.2 | N.D. | U,ICH | 10 | 5.0 | 2.2 | N.D. | U | 10 | 5.0 | 2.2 |
| | Butanone) Methyl isobutyl ketone (4- Methyl-2-Pentanone) | 2,000 | 1300 | N.D. | U | 10 | 5.0 | 4.4 | N.D. | U | 10 | 5.0 | 4.4 | N.D. | U | 10 | 5.0 | 4.4 | N.D. | U | 10 | 5.0 | 4.4 | N.D. | U | 10 | 5.0 | 4.4 |
| | Methyl tert-butyl Ether | 12 | 5 | N.D. | U | 1 | 0.5 | 0.31 | N.D. | U | 1 | 0.5 | 0.31 | N.D. | U | 1 | 0.5 | 0.31 | N.D. | U | 1 | 0.5 | 0.31 | N.D. | U | 1 | 0.5 | 0.31 |
| | Methylene chloride | 4.8 | 9,100 | N.D. | U | 5 | 1.0 | 0.64 | N.D. | U | 5 | 1.0 | 0.64 | N.D. | U | 5 | 1.0 | 0.64 | N.D. | U | 5 | 1.0 | 0.64 | N.D. | U | 5 | 1.0 | 0.64 |
| | Styrene | 100 | 10 | N.D. | U | 11 | 0.5 | 0.17 | N.D. | U | 1 | 0.5 | 0.17 | N.D. | U | 1 | 0.5 | 0.17 | N.D. | U | 1 | 0.5 | 0.17 | N.D. | U | 1 | 0.5 | 0.17 |
| 500 | Tetrachloroethylene | 5 | 170 | N.D. | U | 5 | 0.5 | 0.39 | N.D. | U | 5 | 0.5 | 0.39 | N.D. | U | 5 | 0.5 | 0.39 | N.D. | U | 5 | 0.5 | 0.39 | N.D. | U | 5 | 0.5 | 0.39 |
| | Toluene trans-1,2- Dichloroethylene | 1,000 100 | 40 260 | N.D. | U | 1 1 | 0.5 | 0.24 0.37 | N.D. N.D. | U | 1 1 | 0.5 0.5 | 0.24 | N.D. N.D. | U | 1 1 | 0.5 0.5 | 0.24 | N.D. N.D. | U | 1 | 0.5 | 0.24 | N.D. | U | 1 | 0.5 | 0.24 |
| | Trichloroethylene | 5 | 310 | N.D. | Ü | 1 | 0.5 | 0.37 | N.D. | U | 1 | 0.5 | 0.37 | N.D. | | 1 | 0.5 | 0.37 | N.D. | U | 1 | 0.5 | 0.37 | N.D. | U | 1 | 0.5 | 0.37 |
| | Vinyl chloride | 2 | 3,400 | N.D. | Ü | 1 | 0.5 | 0.3 | N.D. | Ū | 1 | 0.5 | 0.3 | N.D. | Ū | 1 | 0.5 | 0.3 | N.D. | Ü | 1 | 0.5 | 0.3 | N.D. | Ü | 1 | 0.5 | 0.3 |
| | Xylenes | 10,000 | 20 | N.D. | U | 11 | 1.5 | 0.23 | N.D. | U | 1 | 0.5 | 0.23 | 0.32 | J | 1 | 0.5 | 0.23 | N.D. | U | 1 | 0.5 | 0.23 | N.D. | U | 11 | 1.5 | 0.23 |
| EPA 6020 | Dissolved Lead | 15 | 50,000 | | - | - | - | - | 0.0976 | J | 1 | 0.2 | 0.0898 | N.D. | U | 1 | 0.2 | 0.0898 | N.D. | U | 1 | 0.2 | 0.0898 | N.D. | U | 1 | 0.2 | 0.0898 |
| EPA 200.8 | Total Lead | 15 | 50,000 | 0.211 | J | 1.0 | - | 0.0898 | - | - | | - | - | - | - | - | | - | | | - | | _ | - | | - | _ | - |

The data are in micrograms per liter (μ g/L). Shaded values exceeded the DOH EALs. LOD for this analyte exceeds the DOH EAL Not Analyzed

Not Analyzed

This sample was analyzed by EPA Method 200.8 and therefore does not have an LOD

DL Detection Limit or Method Detection Limit (MDL)

DOH EALS

DOH Tier 1 Environmental Action Levels for groundwater where groundwater is a current drinking water source and surface water is greater than 150 meters from the site (DOH, Fall 2011).

EPA Environmental Protection Agency

HD The chromatographic pattern was inconsistent with the profile of the reference fuel standard.

ICH Initial calibration verification recovery is above the control limit for this analyte.

ICJ Initial calibration verification recovery is below the control limit for this analyte.

Analyte was detected at a concentration below the LOQ and above the DL. Reported value is estimated.

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J LOD LOQ N.D. Q TPH-d TPH-g U

Analyte was detected at a concentration below the LOQ a Limit of Detection Limit of Quantitation Not Detected Qualifiers Total Petroleum Hydrocarbons as diesel Total Petroleum Hydrocarbons as gasoline Undetected at DL and is reported as less than the LOD.

Red Hill LTM, 4Q2014 Status Report Inside Tunnel Wells

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SECTION 3 – DATA QUALITY ASSESSMENT

A data quality assessment, which consists of a review of the overall groundwater sample collection and analysis process, was performed in order to determine whether the analytical data generated met the quality objectives for the project. The data quality assessment was performed in accordance with the approved WP/SAP (ESI, 2012). The field quality control program consisted of standardized sample collection and management procedures, and the collection of field duplicate samples, matrix spike samples, and trip blank samples. The laboratory quality assurance program consisted of the use of standard analytical methods and the preparation and analyses of Matrix Spike [MS]/Matrix Spike Duplicate [MSD] samples, surrogate spikes, blanks, and Laboratory Control Samples [LCSs]/Laboratory Control Sample Duplicates [LCSDs].

3.1 DATA VALIDATION AND ASSESSMENT

The objective of data validation is to provide data of known quality for project decisions. Data quality is judged in terms of Precision, Accuracy, Representativeness, Completeness, Comparability, and Sensitivity [PARCCS]. A number of factors may affect the quality of data, including: sample collection methods, sample analysis methods, and adherence to established procedures for sample collection, preservation, management, shipment, and analysis.

Precision

Precision is defined as the reproducibility of replicate measurements. Precision is evaluated by Relative Percentage Difference [RPD] of field duplicates, LCS/LCSD, and MS/MSD results. Field duplicate and MS/MSD samples were collected at a rate of approximately 10% of primary samples. Field duplicates were sent to the laboratory along with the primary samples.

The RPDs of detected analytes for the primary and field duplicate samples (ES114 and ES115) are provided in Table 3.1. A precision of less than 50% for duplicate pairs is required by the DON Project Procedures Manual to be considered acceptable (DON 2007). All duplicate RPDs are less than the acceptable maximum, except for lead. An RPD of 200% was assigned to the ES114 and ES115 (RHMW02) lead results because lead was detected in only one of the replicate samples from the duplicate pair. The lead result for sample ES115 was below the limit of quantitation [LOQ]; consequently, the RPD exceedance signified the anticipated decrease in precision below the LOQ, but was not indicative of a QC issue. Therefore, it is unlikely to have an effect on data usability. In addition, all RPDs for MS/MSD and LCS/LCSD pairs were also within the control limits, and the data precision is considered acceptable.

Accuracy

Accuracy is defined as the degree of conformity of a measurement to a standard or true value. Accuracy is evaluated through measurement of the percent recovery of an analyte in a reference standard or spiked sample. Accuracy limits for surrogates, laboratory control spike, MS, and MSD samples are either prescribed by the Department of Defense [DoD] or

established by the individual laboratory. The acceptance criteria for accuracy are dependent on the analytical method and are based on historical laboratory or DoD data.

Between July 2006 and July 2010, naphthalene was analyzed for by both EPA Methods 8260B and 8270C, and both results were reported. In September 2005 and in all data beginning in October 2010, only results using EPA Method 8270C were reported. Naphthalene has historically only been detected at concentrations above the DOH EALs in well RHMW02. In this well, concentrations of naphthalene detected in each sample by EPA Method 8260B were generally two to three times higher than those detected by EPA Method 8270C. We assume this is due to the better preservation of VOCs associated with the use of EPA Method 8260B. This suggests that the naphthalene results provided by EPA Method 8270C may be biased low. Since March 2014, naphthalene concentrations in RHMW02 have exceeded DOH EALs for both gross contamination and drinking water toxicity. Therefore, a low bias is unlikely to affect project decisions.

Results for TPH-d in samples ES113 through ES118 were flagged "HD." The laboratory indicated a mismatch between the calibration standard and the TPH-d chromatographic profile. Mismatches of this type are not uncommon. Even though chromatograms are not part of the standard laboratory package, ESI was able to review the chromatograms from RHMW02 dating back to October 2012. The chromatograms of groundwater samples from RHMW02 did not significantly differ between each event, but did not match a standard chromatogram of JP-8 in groundwater.

The MS and MSD recoveries were above the control limits for TPH-d, and the associated sample results may be biased high. The TPH-d results from this sample delivery group (ES113 to ES116) have been flagged "J", estimated, to indicate the increased error caused by possible matrix interference. Naphthalene, 2-methylnaphthalene and 1-methylnaphthalene concentrations for ES114, the primary sample on which the MS/MSD were performed, were significantly higher than the added spike concentration, which prevented an accurate evaluation of the MS/MSD recovery for these analytes.

All other MS/MSD recoveries and all LCS and surrogate spike recoveries were within acceptable recovery limits; therefore, the data accuracy for this monitoring event is considered acceptable.

Representativeness

Representativeness is the degree that data accurately and precisely represents a characteristic of a population, parameter variations at a sampling point, or an environmental condition. Representativeness was achieved by conducting sampling in accordance with the sample collection procedures described in the project WP/SAP, including standardized sample collection methods (ESI, 2012).

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Representativeness is also evaluated through the compliance with the standardized sample holding time and sample preservation methods, and through the analysis of blank samples, including method blank and trip blank samples. For this sampling event, all sample holding time and sample preservation were consistent with EPA guidance.

For this sampling event, one trip blank was included in every cooler containing samples for VOC and TPH-g analysis to assess the potential for contamination during sample transport. Two trip blanks were collected. 1,4-Dichlorobenzene was the only COPC detected, and in only one of the trip blanks at a concentration below the LOD; however, 1,4-dichlorobenzene was not detected in any of the groundwater samples, excluding the possibility that project samples were contaminated with 1,4-dichlorobenzene during handling or transport. Based on the assessment of representativeness, the groundwater sample data are considered representative of the groundwater quality on site. The trip blank results are provided in Table 3.1.

Completeness

Completeness is defined as the overall percentage of valid analytical results (including estimated results) compared to the total number of analytical results reported by the analytical laboratory. No data were rejected for this project, and therefore the completeness goal for this project (90%), was successfully met.

Comparability

Comparability expresses the confidence with which one data set can be compared to another data set. Comparability can be related to accuracy and precision because these quantities are measures of data reliability. Data with acceptable precision and accuracy are considered comparable if collection techniques, analytical procedures, methods and reporting are equivalent.

As noted above, between July 2006 and July 2010, naphthalene was analyzed for using both EPA Methods 8260B and 8270C, and in September 2005 and between October 2010 and the most recent event, only results using EPA Method 8270C were reported. In general, EPA Method 8260B resulted in higher, and as discussed above, likely more accurate, results than EPA Method 8270C. However, for the sake of comparability with results from recent events, EPA Method 8270C was used for naphthalene analysis in this event. Consequently, the low bias associated with Method 8270C should be considered when making project decisions.

All project samples for TPH-g analysis through July 2010 were analyzed by EPA Method 8015; beginning in October 2010, EPA Method 8260B was used. There was no event where both methods were used; consequently, there is no way to directly compare the results obtained by method and to assess potential bias. However, there is no reason to believe that using either method should bias the data, and the TPH-g data for all events should be comparable.

Other than the naphthalene bias discussed above, no issues with comparability were identified. The results are considered comparable within this data set and with the data collected from recent sampling events.

Sensitivity

The LOQs are established by the laboratory based on the limits of detection [LODs] or instrument detection limits, historical data, and EPA limits established for the various methods. The LOQs for samples may require adjustment by the laboratory due to matrix interference or if high levels of target analytes necessitate dilution before analysis. Matrix interference and sample dilutions have the effect of decreasing sensitivity and increasing the LOQs. Laboratory LODs and LOQs for several analytes (EPA Methods 8260 and 8270) for this event differed from the LODs and LOQs in the WP/SAP because the laboratory updates them quarterly and in some cases, dilution was necessary due to the presence of high concentrations of analytes.

For this event, LODs and LOQs for several analytes were greater than the DOH EALs (as stated in the WP/SAP), and therefore it would not be possible to detect the analytes at concentrations greater than the DOH EALs but below the LODs. The lack of the required sensitivity should be considered when making project decisions. The affected analytes for this monitoring event are 1,2,3-trichloropropane, 1,2-dibromo-3-chloropropane, 1,2-dibromoethane, 1,2-dichloropropene, bromodichloromethane, chloromethane, dibromochloromethane, 1,1,2,2-tetrachloroethane, and dibenzo[a,h]anthracene.

3.2 DATA ASSESSMENT AND USABILITY CONCLUSIONS

The PARCCS criteria were evaluated, and with a few exceptions, all criteria were met. Based on the high MS/MSD recoveries for TPH-d, it is possible that sample results for TPH-d may be biased high. This increased error should be kept in mind when comparing TPH-d results from this sampling event to those of previous events, but should not significantly affect data usability. Other than this issue, the data assessment concludes that all data generated during this event are usable for the intended purpose.

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TABLE 3.1 Quality Control Results for Groundwater Sampling (October 27 and 28, 2014) Red Hill Bulk Fuel Storage Facility October 2014 Quarterly Monitoring Report

| | | | | 1 | | | | CI EU 17 | Quarteri | | | | | | | | | | | | | 140000 | ~ . | |
|-----------|---|--|------------------------|--------------|--|--|-------------|-------------|--------------|------------------|--------------|-------------|-------------|-----------|--|--|--------------|------------|--------------|--------------|--|---|--|--------------|
| | | DOH EALS RHMW02 (ES114) RHMW02 (ES115) (DUP) | | | | | | | | | | | | | | ES Tri | p (10/27/2 | (014) | | | ES III | ip (10/28/2 | V14) | , , |
| Markena | Chamical Constituent | | | | | | | 1 | | | | | | RPD | | 1 | | | | | 1 | | | |
| Method | Chemical Constituent | Drinking Water Toxicity | Gross Contamination | Result | a | LOQ | LOD | DL | Result | a | LOQ | LOD | DL | Duplicate | Result | Q | LOQ | LOD | DL | Result | l a | LOQ | LOD | DL |
| | | Dimming fraction toxicity | 0.000 00.100.110.100.1 | | | | | | | | | | | (%) | | 1 ~ | | | | itesuit | _ | | | , ' |
| EPA 8015B | TPH-d | 100 | 100 | 2.000 | LID I | 25 | 12 | 11 | 2.000 | LID I | 26 | 12 | 11 | 0.00 | | - | | | | | + | + | - | |
| | L | 190 100 | 100 | 2,000 | HD,J | 25 | 12 | 11 | 2,000 | HD,J | 25 | 12 | 11 | 7.27 | - 10 | - U | - | - 20 | - 20 | - N.D. | | | | - 26 |
| EPA 8260B | TPH-g | 370 | 20 | 0.53 | | 50 0.19 | 0.047 | 26 0.025 | 53 0.53 | + | 50 0.19 | 30 0.047 | 26 0.025 | 0.00 | N.D. | | 50 | 30 | 26 | N.D. | U | 50 | 30 | 26 |
| | Acenaphthene Acenaphthylene | 240 | 2,000 | N.D. | l u | 0.19 | 0.047 | 0.042 | N.D. | u | 0.19 | 0.047 | 0.042 | NA | | + | | | | | | | | |
| | | 1.800 | 2,000 | | U | 0.19 | 0.047 | | N.D. | - | 0.19 | | <u> </u> | NA NA | | + | | | | | - | + | | |
| | Anthracene | 0.092 | 4.7 | N.D. N.D. | U | | | 0.027 | N.D. | U | | 0.047 | 0.027 | NA NA | <u> </u> | - | - | | - | - | | - | ├─ ─┤ | |
| | Benzo[a]anthracene | 1,500 | 0.13 | N.D. | U | 0.19 0.19 | 0.047 | 0.031 | N.D. | U | 0.19 0.19 | 0.047 | 0.031 | NA NA | - | - | - | - | - | - | - | | | <u> </u> |
| | Benzo[g,h,i]perylene | | | | | 0.19 | | | | U | | | | | ļ | + | | - | - | - | - | - | - | - |
| | Benzo[a]pyrene | 0.2 | 0.81 0.75 | N.D. | U | 0.19 | 0.047 | 0.021 | N.D. | l U | 0.19 | 0.047 | 0.021 | NA NA | - | - | - | | - | - | | + | | |
| | Benzo[b]fluoranthene | 0.92 | 0.75 | N.D. | T U | 0.19 | 0.047 | 0.017 | N.D. | U | 0.19 | 0.047 | 0.017 | NA NA | | - | | - | - | - | | - | - + | |
| | Benzo[k]fluoranthene | 9.2 | 1 | N.D. | 1 0 | 0.19 | 0.047 | 0.029 | N.D. | U | 0.19 | 0.047 | 0.029 | NA NA | - | + | | - | - | - | | - | | <u> </u> |
| EPA 8270C | Chrysene Dibenzo[a,h]anthracene | 0.0092 | 0.52 | N.D. | U | 0.19 | 0.047 | 0.024 | N.D. | U | 0.19 | 0.047 | 0.025 | NA NA | | + | | - | - | | | | | |
| | Fluoranthene | 1,500 | 130 | N.D. | | 0.19 | 0.047 | 0.043 | N.D. | 1 0 | 0.19 | 0.047 | 0.043 | NA NA | - | | - | | - | | | - | | <u> </u> |
| | Fluorene | 240 | 950 | N.D. | 1 0 | 0.19 | 0.047 | 0.044 | N.D. | U | 0.19 | 0.047 | 0.044 | NA NA | - | + | | - | - | - | | - | | <u> </u> |
| | | 0.092 | 0.095 | N.D. | | 0.19 | 0.047 | 0.020 | N.D. | ŭ | 0.19 | 0.047 | 0.020 | NA NA | - | + | | - | - | | | - | | - |
| | Indeno[1,2,3-cd]pyrene 1-Methylnaphthalene | 4.7 | 10 | N.D. | | 3.8 | 1.9 | 0.020 | N.D. 54 | 1- | 3.8 | 1.9 | 0.020 | 8.85 | | - | - | - | - | - | + | + | - | <u> </u> |
| | 2-Methylnaphthalene | 24 | 10 | 43 | | 3.8 | 0.95 | 0.88 | 36 | 1 | 3.8 | 0.95 | 0.98 | 17.72 | | + | | - | - | - | | + | | - |
| | Naphthalene | 17 | 21 | 140 | - | 3.8 | 0.95 | 0.64 | 130 | - | 3.8 | 0.95 | 0.64 | 7.41 | | - | - | - | - | - | - | - | + | <u> </u> |
| | Phenanthrene | 240 | 410 | N.D. | U | 0.19 | 0.95 | 0.026 | N.D. | U | 0.19 | 0.95 | 0.026 | NA NA | - | - | | - | - | - | | | | - |
| | Pyrene | 180 | 68 | N.D. | T U | 0.19 | 0.047 | 0.028 | N.D. | U | 0.19 | 0.047 | 0.028 | NA NA | | + | | - | - | - | - - | | 1 | - |
| | 1,1,1,2-Tetrachloroethane | 0.52 | 50,000 | N.D. | T U | 1 | 0.5 | 0.4 | N.D. | 1 0 | 1 | 0.5 | 0.4 | NA NA | N.D. | T U | 1 | 0.5 | 0.4 | N.D. | | + | 0.5 | 0.4 |
| | 1,1,2,2-Tetrachloroethane | 0.067 | 500 | N.D. | T U | | 0.5 | 0.41 | N.D. | T U | 1 | 0.5 | 0.41 | NA NA | N.D. | T U | 1 | 0.5 | 0.41 | N.D. | l ü | | 0.5 | 0.41 |
| | 1,1,1-Trichloroethane | 200 | 970 | N.D. | T Ü | 5 | 0.5 | 0.3 | N.D. | l ü l | 5 | 0.5 | 0.3 | NA NA | N.D. | T U | <u>'</u> | 0.5 | 0.3 | N.D. | T U | 5 | 0.5 | 0.3 |
| | 1,1,2-Trichloroethane | 5 | 50,000 | N.D. | T U | 1 | 0.5 | 0.38 | N.D. | l ŭ l | 1 | 0.5 | 0.38 | NA NA | N.D. | T Ü | 1 | 0.5 | 0.38 | N.D. | l ü | 1 | 0.5 | 0.38 |
| | 1.1-Dichloroethane | 2.4 | 50,000 | N.D. | U | 5 | 0.5 | 0.28 | N.D. | T U | 5 | 0.5 | 0.28 | NA NA | N.D. | TÜ | <u>-</u> | 0.5 | 0.28 | N.D. | T Ü | 5 | 0.5 | 0.28 |
| | 1,1-Dichloroethylene | 7 | 1,500 | N.D. | Ü | 1 | 0.5 | 0.43 | N.D. | T Ü | 1 | 0.5 | 0.43 | NA | N.D. | T U | 1 | 0.5 | 0.43 | N.D. | l ŭ | 1 | 0.5 | 0.43 |
| | 1,2,3-Trichloropropane | 0.6 | 50.000 | N.D. | T Ū | 5 | 1 | 0.64 | N.D. | T Ü | 5 | 1 | 0.64 | NA | N.D. | Ü | <u>·</u> | 1 | 0.64 | N.D. | l ü | 5 | 1 | 0.64 |
| | 1,2,4-Trichlorobenzene | 70 | 3,000 | N.D. | T Ū | 5 | 1 | 0.5 | N.D. | T Ü | 5 | 1 | 0.5 | NA | N.D. | T U | 5 | 1 | 0.5 | N.D. | Ť | 5 | 1 1 | 0.5 |
| | 1,2-Dibromo-3- chloropropane | 0.04 | 10 | N.D. | T Ū | 10 | 2 | 1.2 | N.D. | l ü l | 10 | 2 | 1.2 | NA | N.D. | t ü | 10 | 2 | 1.2 | N.D. | T Ū | 10 | 2 | 1.2 |
| | 1,2-Dibromoethane | 0.04 | 50,000 | N.D. | Ū | 1 | 0.5 | 0.24 | N.D. | Ū | 1 | 0.5 | 0.24 | NA | N.D. | Ū | 1 | 0.5 | 0.24 | N.D. | Ū | 1 | 0.5 | 0.24 |
| | 1,2-Dichlorobenzene | 600 | 10 | N.D. | Ū | 1 | 0.5 | 0.46 | N.D. | l ū l | 1 | 0.5 | 0.46 | NA | N.D. | Ū | 1 | 0.5 | 0.46 | N.D. | l ū | | 0.5 | 0.46 |
| | 1,2-Dichloroethane | 0.15 | 7,000 | N.D. | U,IH | 1 | 0.5 | 0.24 | N.D. | U,IH | 1 | 0.5 | 0.24 | NA | N.D. | U,IH | 1 | 0.5 | 0.24 | N.D. | U | 1 | 0.5 | 0.24 |
| | 1,2-Dichloropropane | 5 | 10 | N.D. | Ü | 5 | 0.5 | 0.42 | N.D. | Ú | 5 | 0.5 | 0.42 | NA | N.D. | Ú | 5 | 0.5 | 0.42 | N.D. | U | 5 | 0.5 | 0.42 |
| | 1,3-Dichlorobenzene | 180 | 5 | N.D. | U | 1 | 0.5 | 0.4 | N.D. | T U | 1 | 0.5 | 0.4 | NA | N.D. | U | 1 | 0.5 | 0.4 | N.D. | U | 1 | 0.5 | 0.4 |
| | 1,3-Dichloropropene (total of cis/trans) | 0.43 | 50,000 | N.D. | U | 1 | 0.5 | 0.25 | N.D. | U | 1 | 0.5 | 0.25 | NA | N.D. | U | 1 | 0.5 | 0.25 | N.D. | U | 1 | 0.5 | 0.25 |
| | 1,4-Dichlorobenzene | 75 | 5 | N.D. | U | 1 | 0.5 | 0.43 | N.D. | U | 1 | 0.5 | 0.43 | NA | N.D. | U | 1 | 0.5 | 0.43 | 0.43 | U | 1 | 0.5 | 0.43 |
| | Acetone | 22,000 | 20,000 | N.D. | U,ICH | 20 | 10 | 6 | N.D. | U,IC | 20 | 10 | 6 | NA | N.D. | U,IC | 20 | 10 | 6 | N.D. | U,IC | 20 | 10 | 6 |
| | Benzene | 5 | 170 | N.D. | U | 1 | 0.5 | 0.14 | N.D. | U | 1 | 0.5 | 0.14 | NA | N.D. | U | 1 | 0.5 | 0.14 | N.D. | U | 1 | 0.5 | 0.14 |
| | Bromodichloromethane | 0.12 | 50,000 | N.D. | U | 5 | 0.5 | 0.21 | N.D. | U | 5 | 0.5 | 0.21 | NA | N.D. | U | 5 | 0.5 | 0.21 | N.D. | U | 5 | 0.5 | 0.21 |
| | Bromoform | 80 | 510 | N.D. | U | 10 | 1 | 0.5 | N.D. | U | 10 | 1 | 0.5 | NA | N.D. | U | 10 | 1 | 0.5 | N.D. | U | 10 | 1 | 0.5 |
| EPA 8260B | Bromomethane | 8.7 | 50,000 | N.D. | U | 20 | 5 | 3.9 | N.D. | U | 20 | 5 | 3.9 | NA | N.D. | U | 20 | 5 | 3.9 | N.D. | U,ICJ | 20 | 5 | 3.9 |
| | Carbon Tetrachloride | 5 | 520 | N.D. | U | 1 | 0.5 | 0.23 | N.D. | U | 1 | 0.5 | 0.23 | NA | N.D. | U | 1 | 0.5 | 0.23 | N.D. | U | 1 | 0.5 | 0.23 |
| | Chlorobenzene | 100 | 50 | N.D. | U | 5 | 0.5 | 0.17 | N.D. | U | 5 | 0.5 | 0.17 | NA | N.D. | U | 5 | 0.5 | 0.17 | N.D. | U | 5 | 0.5 | 0.17 |
| | Chloroethane | 21,000 | 16 | N.D. | U | 10 | 5 | 2.3 | N.D. | U | 10 | 5 | 2.3 | NA | N.D. | U | 10 | 5 | 2.3 | N.D. | U | 10 | 5 | 2.3 |
| | Chloroform | 70 | 2,400 | N.D. | U | 5 | 0.5 | 0.46 | N.D. | U | 5 | 0.5 | 0.46 | NA | N.D. | U | 5 | 0.5 | 0.46 | N.D. | U | 5 | 0.5 | 0.46 |
| | Chloromethane | 1.8 | 50,000 | N.D. | U, IJ | 10 | 2 | 1.8 | N.D. | U, IJ | 10 | 2 | 1.8 | NA | N.D. | U, IJ | 10 | 2 | 1.8 | N.D. | U | 10 | 2 | 1.8 |
| | cis-1,2-Dichloroethylene | 70 | 50,000 | N.D. | U | 1 | 0.5 | 0.48 | N.D. | U | 1 | 0.5 | 0.48 | NA | N.D. | U | 11 | 0.5 | 0.48 | N.D. | U | 1 | 0.5 | 0.48 |
| | Dibromochloromethane | 0.16 | 50,000 | N.D. | U | 1 | 0.5 | 0.25 | N.D. | U | 1 | 0.5 | 0.25 | NA NA | N.D. | U | 1 | 0.5 | 0.25 | N.D. | U | 1 | 0.5 | 0.25 |
| | Ethylbenzene | 700 | 30 | 0.15 | J | 1 1 | 0.5 | 0.14 | 0.14 | | 1 | 0.5 | 0.14 | 6.90 | N.D. | U | 1 | 0.5 | 0.14 | N.D. | U | 1 1 | 0.5 | 0.14 |
| | Hexachlorobutadiene | 0.86 | 6 | N.D. | U | 1 1 | 0.5 | 0.32 | N.D. | U | 1 | 0.5 | 0.32 | NA NA | N.D. | U | 1 10 | 0.5 | 0.32 | N.D. | U | 1 1 | 0.5 | 0.32 |
| | Methyl ethyl ketone (2-Butanone) | 7,100 | 8,400 | N.D. | U,ICH | 10 | 5.0 | 2.2 | N.D. | U,IC | 10 | 5.0 | 2.2 | NA NA | N.D. | U,IC | 10 | 5.0 | 2.2 | N.D. | U | 10 | 5.0 | 2.2 |
| | Methyl isobutyl ketone (4-Methyl-2-Pentanone) | 2,000 | 1300 | N.D. | U | 10 | 5.0 | 4.4 | N.D. | U | 10 | 5.0 | 4.4 | NA NA | N.D. | U. | 10 | 5.0 | 4.4 | N.D. | U | 10 | 5.0 | 4.4 |
| | Methyl tert-butyl Ether | 12 | 5 | N.D. | U | 1 | 0.5 | 0.31 | N.D. | U | 1 | 0.5 | 0.31 | NA NA | N.D. | U | 1 | 0.5 | 0.31 | N.D. | U | | 0.5 | 0.31 |
| | Methylene chloride | 4.8 | 9,100 | N.D. | U | 5 | 1.0 | 0.64 | N.D. | l U | 5 | 1.0 | 0.64 | NA NA | N.D. | U | 5 | 1.0 | 0.64 | N.D. | l U | 5 | 1.0 | 0.64 |
| | Styrene | | 10 170 | N.D. | U | 1 6 | 0.5 | 0.17 | N.D. | U | 1 6 | 0.5 | 0.17 | NA NA | N.D. | l U | 1 | 0.5 | 0.17 | N.D. | U | 1 5 | 0.5 | 0.17 |
| | Tetrachloroethylene Toluene | 5 1,000 | 40 | N.D. | U | 5 | 0.5 | 0.39 | N.D. N.D. | U | 5 1 | 0.5 | 0.39 | NA NA | N.D. N.D. | U | <u>5</u> | 0.5 0.5 | 0.39 0.24 | N.D. N.D. | U | 5 | 0.5 | 0.39 |
| | trans-1,2- Dichloroethylene | 1,000 | 260 | N.D. | U | 1 | 0.5 | 0.24 | N.D. | U | 1 | 0.5 | 0.24 | NA NA | N.D. | U | 1 | 0.5 | 0.24 | N.D. | U | + | 0.5 | 0.24 |
| | Trichloroethylene | 5 | 310 | N.D. | T U | 1 | 0.5 | 0.37 | N.D. | U | 1 | 0.5 | 0.37 | NA NA | N.D. | 1 0 | 1 | 0.5 | 0.37 | N.D. | U | 1 | 0.5 | 0.37 |
| | Vinyl chloride | 2 | 3,400 | N.D. | T U | 1 | 0.5 | 0.37 | N.D. | l ü l | 1 | 0.5 | 0.37 | NA NA | N.D. | U | 1 | 0.5 | 0.37 | N.D. | U | | 0.5 | 0.37 |
| | Xylenes | 10,000 | 20 | 0.32 | | 1 | 0.5 | 0.23 | 0.29 | J | 1 | 0.5 | 0.23 | 9.84 | N.D. | 1 0 | 11 | 1.5 | 0.23 | N.D. | | 11 | 1.5 | 0.23 |
| | | · · · · · · · · · · · · · · · · · · · | 50.000 | N.D. | l U | | 0.3 | 0.0898 | 0.165 | + | | | 0.0898 | | | + | | 1.5 | 0.20 | | + | + | + ' | 0.23 |
| EPA 6020 | Lead | 15 | 50 000 | | | | | | | | 1 | 0.2 | ()()898 | 200 | | | - | | - | - | | | | |

The data are in micrograms per liter (µg/L). Shaded values exceeded the DOH EALs.

Not Analyzed

DOH Tier 1 Environmental Action Levels for groundwater where groundwater is a current drinking water source and surface water is greater than 150 meters from the site (DOH, Fall 2011).

DL Detection Limit or Method Detection Limit (MDL)

EPA Environmental Protection Agency

HD The chromatographic pattern was inconsistent with the profile of the reference fuel standard.

ICH Initial calibration verification recovery is above the control limit for this analyte.

ICJ Initial calibration verification recovery is below the control limit for this analyte.

J Analyte was detected at a concentration below the LOQ and above the DL. Reported value is estimated.

LOD LOQ NA N.D. Q TPH-g TPH-d

Limit of Detection
Limit of Quantitation
Both results for duplicate pair were non-detect, no RPD calculations
Not Detected
Qualifiers
Total Petroleum Hydrocarbons as gasoline
Total Petroleum Hydrocarbons as diesel
Undetected at DL and is reported as less than the LOD.

3-5

January 2015

Contract No. N62742-12-D-1853

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SECTION 4 - SUMMARY, CONCLUSIONS, AND RECOMMENDATIONS

On October 27 and 28, 2014, ESI personnel collected groundwater samples from four monitoring wells at the RHSF (wells RHMW01, RHMW02, RHMW03, and RHMW05) and one sampling point at Red Hill Shaft (RHMW2254-01).

The groundwater sampling was conducted as part of the long-term groundwater and soil vapor monitoring program at the RHSF, under NAVFAC Contract Number N62742-12-D-1853. The sampling was conducted in accordance with the approved WP/SAP prepared by ESI. A summary of the analytical results is provided below.

- RHMW01 TPH-d (120 µg/L) and lead (0.0976 µg/L) were the only analytes detected. The
 TPH-d concentration exceeded the DOH EAL for gross contamination, but was below the
 SSRBL of 4,500 µg/L for TPH. None of the other chemical constituents analyzed for were
 detected at concentrations above their respective DOH EAL.
- RHMW02 TPH-d (2,000 μg/L), TPH-g (57 and 53 μg/L), xylenes (0.32 and 0.29 μg/L), acenaphthene (0.53 μg/L), ethylbenzene (0.15 and 0.14 μg/L), 1-methylnaphthalene (59 and 54 μg/L), 2-methylnaphthalene (43 and 36 μg/L), and naphthalene (140 and 130 μg/L) were detected in both the primary and duplicate samples collected. Lead (0.165 μg/L) was only detected in the duplicate sample. TPH-d, 1-methylnaphthalene, 2-methylnaphthalene, and naphthalene were detected at concentrations above their respective DOH EALs for both drinking water toxicity and gross contamination. However, the TPH-d concentrations did not exceed the SSRBL.
- RHMW03 TPH-d (80 μg/L) was the only analyte detected. The concentration did not exceed the DOH EAL or the SSRBL.
- RHMW05 TPH-d (16 μ g/L) was the only analyte detected. The concentration did not exceed the DOH EAL or the SSRBL.
- RHMW2254-01 TPH-d (22 μg/L) and total lead (0.211 μg/L) were the only analytes detected. The concentrations did not exceed their respective DOH EAL.

Groundwater Contaminant Trends

- RHMW01 COPCs detected during this round of quarterly sampling are consistent with the
 historical data for RHMW01. TPH-d has historically been detected at concentrations above
 the DOH EAL for both drinking water toxicity and gross contamination. TPH-d
 concentrations have shown an overall decreasing trend from a high of 1,500 µg/L in
 February 2005. The TPH-d concentration exceeded the DOH EAL for gross contamination
 for the first time since April 7, 2014.
- RHMW02 COPCs detected during this round of quarterly sampling are consistent with the
 historical data for RHMW02. TPH-g, TPH-d, 1-methylnaphthalene, 2-methylnaphthalene,
 and naphthalene have historically been detected at concentrations above the DOH EALs.
 During the October 2014 event, concentrations of TPH-d, 1-methylnaphthalene,

2-methylnaphthalene, and naphthalene approximately doubled from the previous event in July 2014 and remained above DOH EALs. The concentrations of 1-methylnaphthalene, 2-methylnaphthalene, and naphthalene increased to their highest levels in over 5 years. The concentrations of these PAHs are also significantly higher than they were in January 2014, when a high TPH-d spike was observed in this well. The concentrations of TPH-g remained below the DOH EALs for gross contamination and drinking water toxicity and were comparable to the concentrations detected during the previous event.

- RHMW03 COPCs detected during this round of quarterly sampling are consistent with the historical data for RHMW03. TPH-d has historically been detected at concentrations above the DOH EALs; however, it has not been detected at concentrations above the DOH EALs since October 2010.
- RHMW05 COPCs detected during this round of quarterly sampling are consistent with the
 historical data for RHMW05. TPH-d has historically been detected in RHMW05 at
 concentrations above the DOH EAL for both drinking water toxicity and gross contamination;
 however, it has not been detected at concentrations above the DOH EALs since January
 2010.
- RHMW2254-01 COPCs detected during this round of quarterly sampling are consistent
 with the historical data for RHMW2254-01. Although the method reporting limits for TPH-d
 exceeded one or both DOH EALs for drinking water toxicity and gross contamination
 between May 2009 and July 2010, TPH-d was last detected in RHMW2254-01 at a
 concentration above the DOH EAL for gross contamination in January 2008.

Conclusions and Recommendations

During the sampling event conducted on October 27 and 28, 2014, TPH-d in RHMW01 and TPH-d, 1-methylnaphthalene, 2-methylnaphthalene, and naphthalene in RHMW02 were detected at concentrations exceeding the DOH EALs. Groundwater contaminant concentrations in wells RHMW03, RHMW05, and RHMW2254-01 remained at low concentrations and did not change significantly from the previous event, or were not detected.

Concentrations of 1-methylnaphthalene, 2-methylnaphthalene, and naphthalene in RHMW02 have shown an increasing trend since March 5, 2014, but have remained below the historic maximums detected in the well. All other analytical results were generally consistent with historical data.

Based on the groundwater monitoring results and the reported release at Tank 5 in January 2014, continued groundwater monitoring at the wells inside the RHSF tunnel is recommended. The next quarterly event is tentatively scheduled for January 2015.

SECTION 5 – FUTURE WORK

Future work includes the first quarter 2015 groundwater monitoring which is tentatively scheduled for January 2015. A quarterly groundwater monitoring report will be prepared to document the sampling event.

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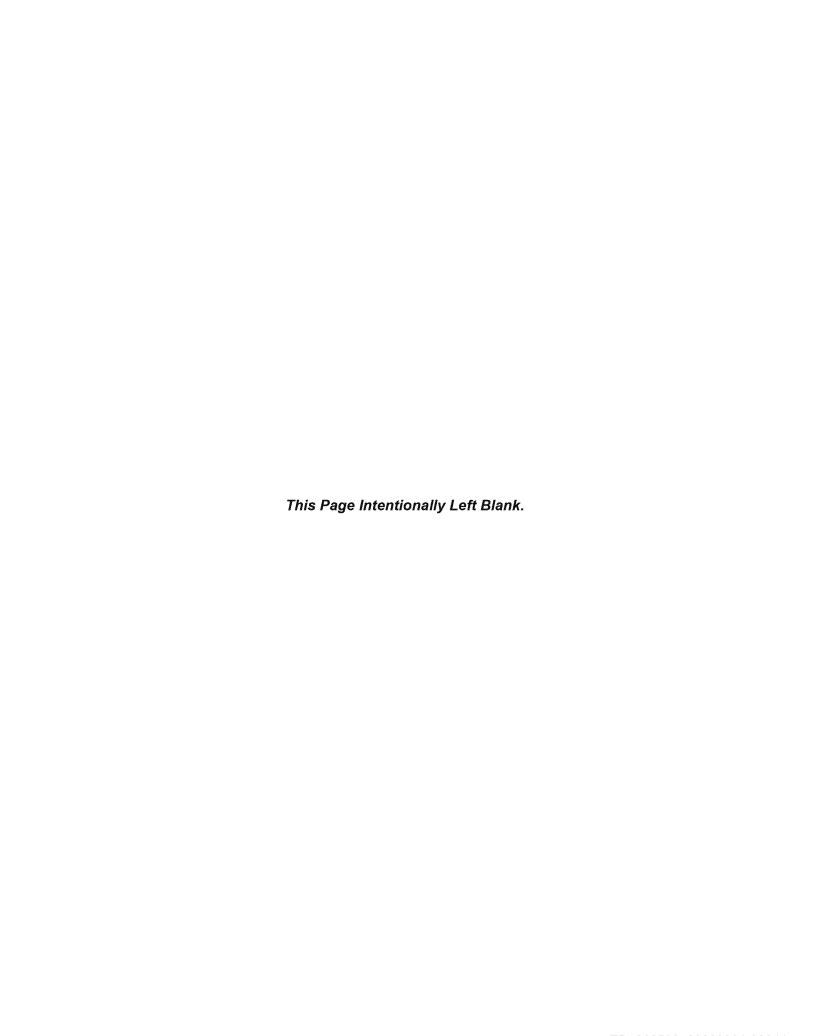
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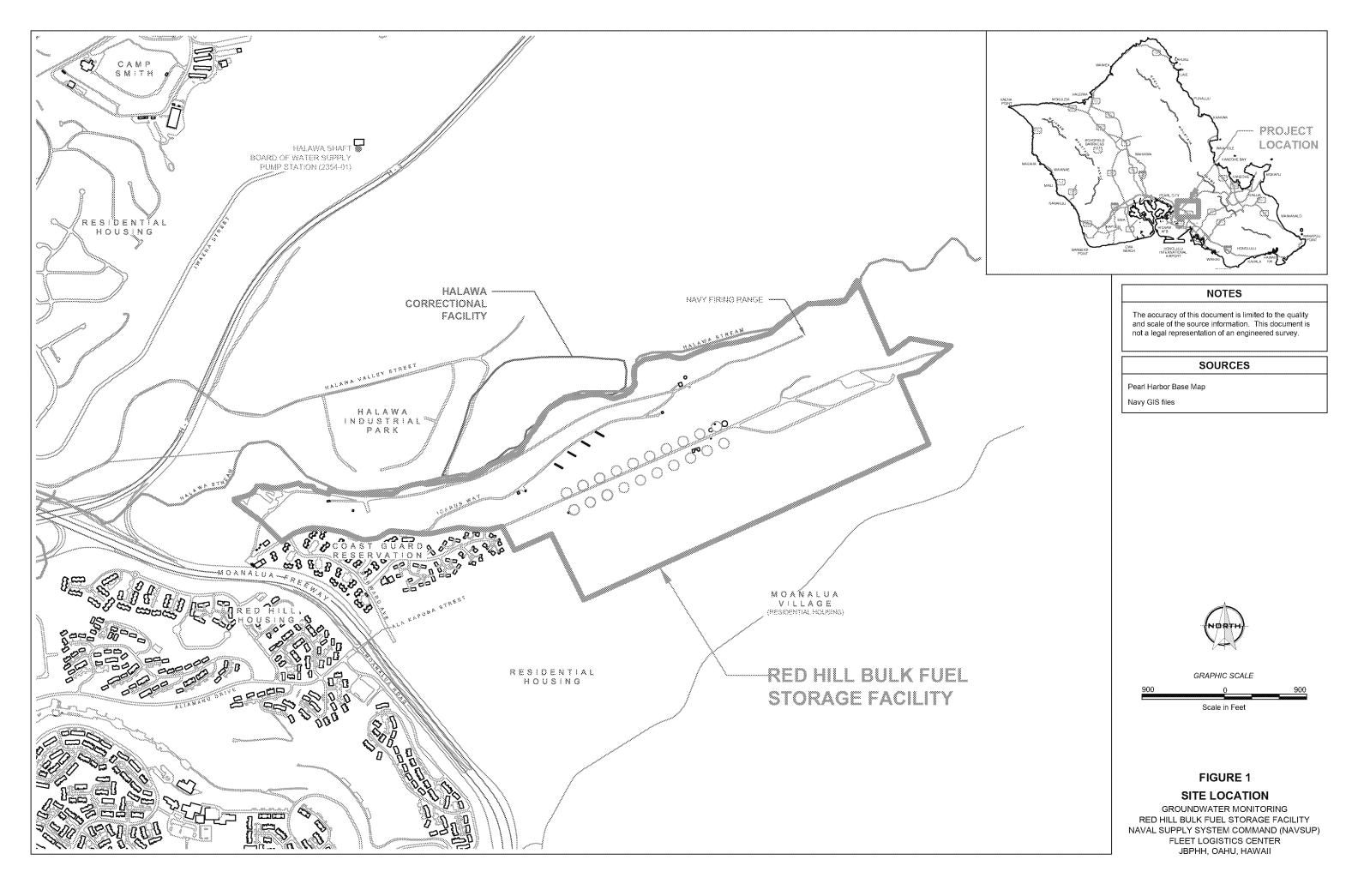
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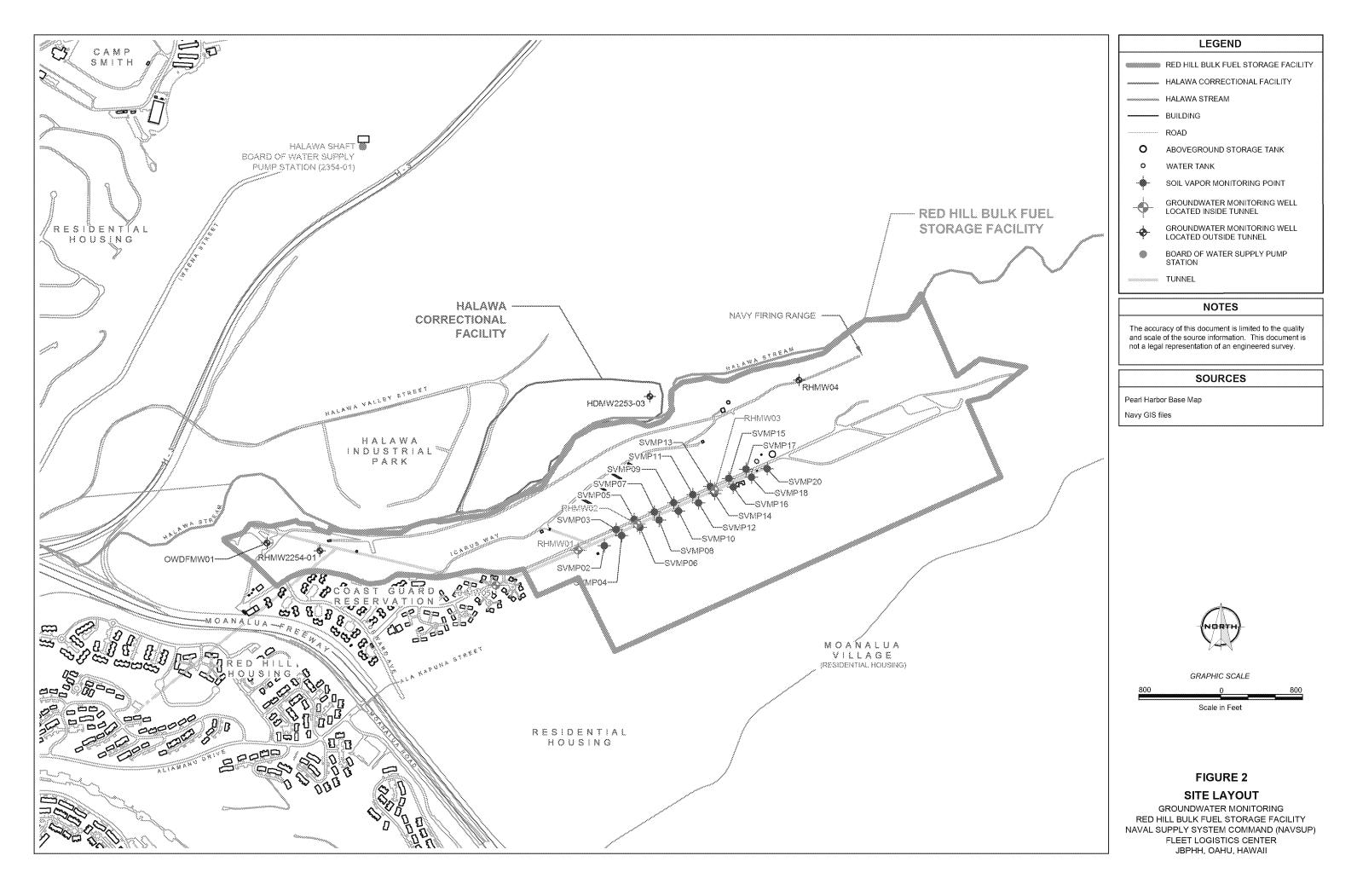
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FIGURES



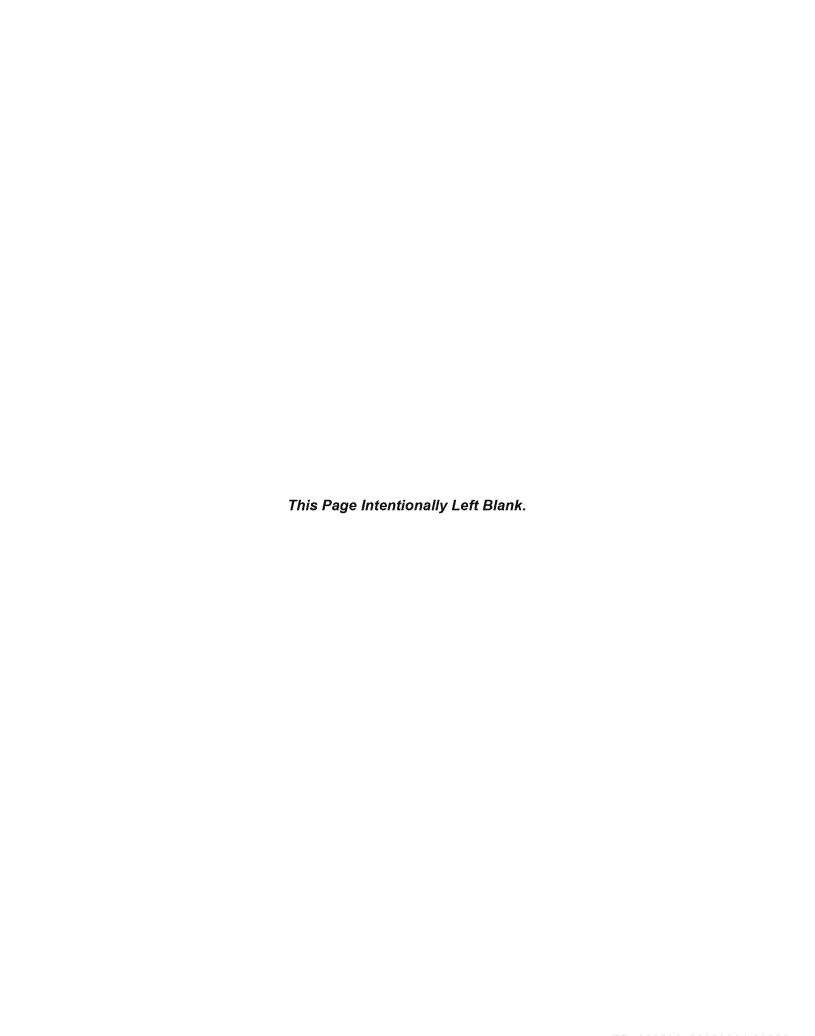






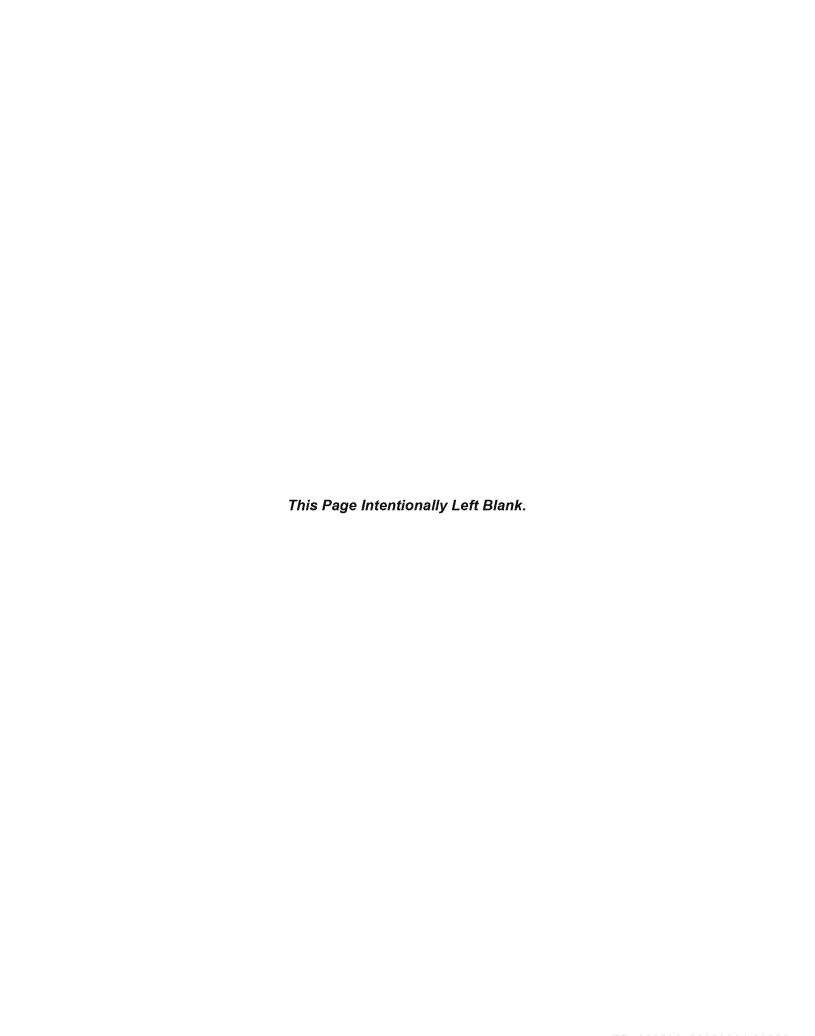


APPENDIX A Groundwater Sampling Logs





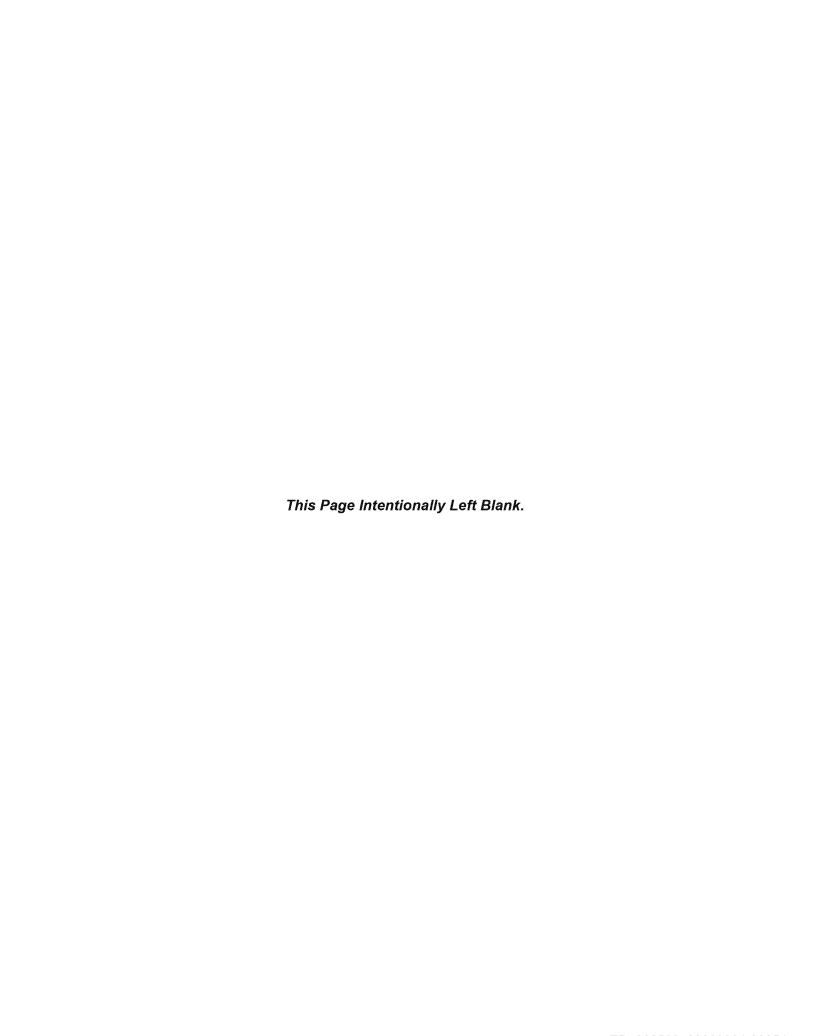
| Well ID: | RHMW01 | | Location: | Red Hil | l Bulk Fuel Stora | ige Facility | Project No.: | 112066 | |
|------------|-----------------------------------|--------------------|-------------------|------------------------|---|----------------------------|-----------------|---------------------|--|
| Initial Wa | Water Level: 83.79 ft | | Date: | 10/27/2014 | | Time:13 | 300 | | |
| Total De | Total Depth of Well: 97.35 ft | | | Personnel Involved: | | Kirk Markle, Jeff Hattemer | | | |
| Length o | ength of Saturated Zone: 13.56 ft | | | Weath | er Conditions: | Not applicable | – well is locat | ted indoors | |
| Volume | of Water to b | e Removed | l: <u>3.5 L</u> | Method | l of Removal: | Bladder Pump | | | |
| Water Le | evel After Pu | ging: | 83.79 ft | Pumpir | ng Rate: | 0.19 L/min | | | |
| Well Pur | ge Data: | | | | | | | | |
| Time | Volum Remov | | | onductivity (mS/cm) | DO (mg/l) | Temperature | Salinity | Redox (ORP) (mV) | |
| 1315 | 0.0 L | 8. | 40 | 0.342 | 11.41 | 25.25 | | 247.8 | |
| 1318 | 0.5 L | 8.3 | 39 | 0.340 | 2.43 | 24.88 | | 289.0 | |
| 1322 | 1.0 L | 8.3 | 38 | 0.339 | 1.64 | 24.23 | _ | -312.0 | |
| 1325 | 1.5 L | 8.3 | 38 | 0.335 | 0.34 | 23.55 | | -346.0 | |
| 1327 | 2.0 L | 8.3 | 37 | 0.335 | -0.07 | 23.41 | | -345.0 | |
| 1329 | 2.5 L | 8.3 | 35 | 0.334 | -0.06 | 23.37 | | -342.1 | |
| 1331 | 3.0 L | 8.3 | 36 | 0.334 | -0.07 | 23.36 | | -341.2 | |
| 1333 | 3.5 L | | 36 | 0.332 | -0.07 | 23.34 | MAX. | -339.6 | |
| | | | | | | | | | |
| | | | | Bladder Pu | mp | | | | |
| ' ' | Color: | | | Clear | | | | | |
| | Turbidit | /: | | Low | | | | | |
| | Sedime | ****************** | | None | | | | | |
| | Other: | | | None | | | | | |
| Laborato | ory Analysis F | arameters' | and Pres | ervatives: | TPH-d - 8015; lead - 6020 | TPH-g, VOCs - 8 | 8260; PAHs - | 8270c sim; | |
| Number | and Types o | f Sample C | ontainers: | 6 - 40m | nl VOAs, 2 - 1L a | ımber jar, 1 - 500 | ıml amber jar, | 1 - 250ml HDPE | |
| Sample | Identification | Numbers: | ES113 | [0400] | | | | | |
| Deconta | mination Pro | cedures: | Triple Rin | sed | | | | | |
| Notes: | YSI did not h | ave salinity | paramete | er. | | | | | |
| Sampled | | Markle, Jeff | Hatteme | ? | | | | | |
| | Delivered to | : Cal | science E | nvironment | *************************************** | Transporters: <u>Fe</u> | dEx | | |
| Date: | 10/27/2014 | | | | | Time: 1530 | | | |
| | | | | - | sing (Gallons/Lii 8"-2.61 • 10"-4. | · | | | |





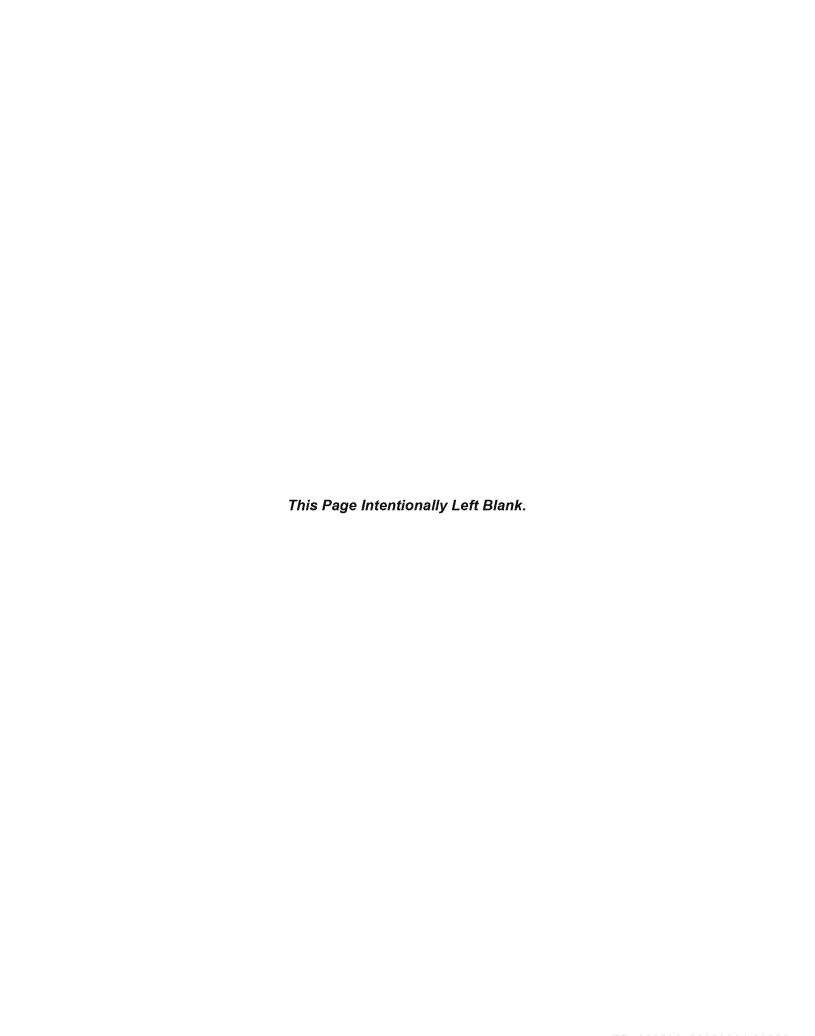
| Well ID | :RHMW02 | | Location: | Red Hil | l Bulk Fuel Stora | ge Facility | Project No.: | 112066 | |
|-------------------------------|--------------------------------------|--|---|---|--|--|---|---------------------|--|
| Initial W | /ater Level: _ | 86.5 | 1 ft | Date: | 10/27/2014 | | Time:1 | 011 | |
| Total Depth of Well: 92.91 ft | | | Person | nel Involved: | Kirk Markle, Jeff Hattemer | | | | |
| Length | of Saturated | Zone: | 6.40 ft | Weath | er Conditions: | Not applicable – well is located indoors | | | |
| Volume | Volume of Water to be Removed: 4.0 L | | | Method | l of Removal: | Bladder Pump | | | |
| Water L | _evel After Pu | ırging: | 86.45 ft | Pumpir | ng Rate: | 0. | 17 L/min | | |
| Well Pu | ırge Data: | | | | | | | | |
| Time | Volur Remo | | рH | Conductivity (mS/cm) | DO (mg/l) | Temperature | Salinity | Redox (ORP) (mV) | |
| 1029 | 9 0.0 L | | 8.40 — | 0.373 | 2.40 | 25.17 | - | | |
| 1035 | | | 8.41 | 0.575 | 1.30 | 24.26 | - | | |
| 1040 | | | 8.40 | 0.551 | 0.97 | 24.13 | | -352.4 | |
| 1045 | | | 8.39 | 0.560 | 0.82 | 24.21 | | -349.7 | |
| 1053 | | | 8.35 | 0.567 | 0.84 | 24.25 | | -361.3 | |
| | | | | | | | | | |
| | | | | | | | | | |
| | | 10000000000000000000000000000000000000 | | *************************************** | MARAGRARARARARARARARARARARARARARARARARAR | 50000000000000000000000000000000000000 | *************************************** | | |
| | | | | | | | | | |
| | | | | | | | | | |
| | | | | | | | | | |
| Sample | Withdrawal I | Method: | : | Bladder Pur | | | | | |
| | ance of Samp | | | | | | | | |
| | Color: | | | Clear | | | | | |
| | Turbidi | ty: | | Low | | | | | |
| | Sedime | | *************************************** | None | ###################################### | | | | |
| | Other: | | | None | | | | | |
| l ahorat | tory Analysis I | Parame | eters and Pres | envatives: | TPH_d _ 8015: | TPH-g, VOCs - | 8260: PAHs - | 8270c sim: | |
| Laborat | ory Arialysis | aramo | | civatives. | lead - 6020 | 11 11-g, v 0 0 3 - | 0200,174113 | 02700 31111, | |
| Numbe | r and Types o | of Samo | le Containers | : 16 - 40ml | VOAs, 6 - 1L am | nber iar. 4 - 500n | nl amber iar. 4 | - 500ml HDPE | |
| | dentification | - | | *********************** | 114 MS/MSD [10 | | | | |
| • | amination Pro | | | | | , (| -/[] | | |
| | | | linity paramet | | | | | | |
| Sample | | | | - | | | | | |
| | ed Delivered to | 0: | Calscience l | Environmenta | al Lab . | Transporters: Fe | edEx | | |
| Date: | 10/27/2014 | | | | | Time: 1530 | | | |
| | | | Ca | apacity of Cas | sing (Gallons/Lin | ear Feet) | | - | |

Capacity of Casing (Gallons/Linear Feet)
2"-0.16• 4"-0.65 • 8"-2.61 • 10"-4.08 • 12"-5.87



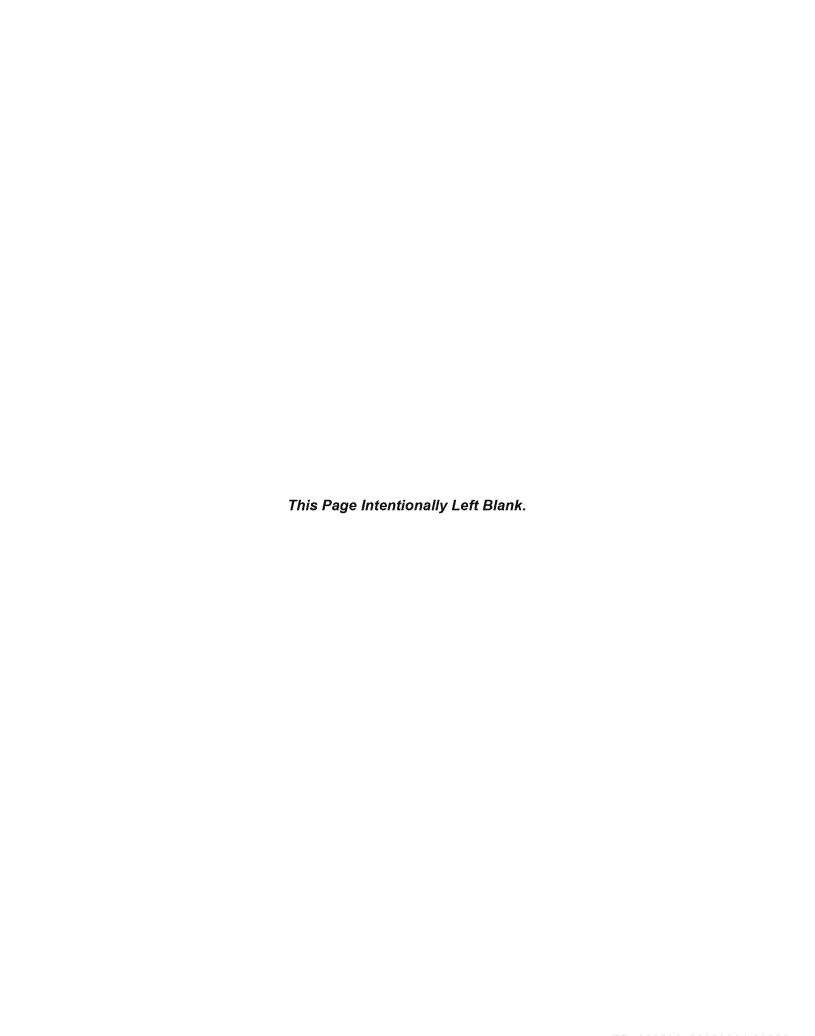


| Well ID: | RHMW03 | | Location: | Red Hil | l Bulk Fuel Stora | age Facility | Project No | o.: <u>1</u> | 12066 |
|--------------------------------------|-----------------------------|---|-------------------------|---|---------------------------------------|---------------------------------|----------------|--------------------|---------------------|
| Initial Wa | itial Water Level:102.78 ft | | | Date: | 10/27/2014 | | Time: | 820 | |
| Total De | pth of Well: | 110 | 0.12 ft | _ Person | nel Involved: | Kirk Markle, Jeff Hattemer | | | |
| Length of Saturated Zone: 7.34 ft | | | | _ Weath | er Conditions: | Not applicabl | e – well is lo | ocated i | ndoors |
| Volume | of Water to b | oe Remove | ed: <u>3.0 L</u> | Method | l of Removal: | Bladder Pump | | | |
| Water Level After Purging: 102.88 ft | | | | _ Pumpir | ng Rate: | C | 0.2 L/min | | |
| Well Pur | • | | | | | | | | |
| Time | Volun Remov | | С <u>рН</u> | onductivity (mS/cm) | DO (mg/l) | Temperature | s Sali | nity | Redox (ORP) (mV) |
| 830 | 0.0 | <u> </u> | 7.52 | 0.823 | 2.14 | 26.10 | | | -206.0 |
| 834 | 1.0 | <u> </u> | ⁷ .56 | 0.820 | 0.95 | 26.40 | | | -240.1 |
| 840 | 2.0 | L 7 | 7.63 | 0.821 | 0.87 | 26.40 | - | | -233.1 |
| 845 | 3.0 | | 7.58 <u> </u> | 0.816 | 0.71 | 26.41 | | | -238.6 |
| | | | | 000000000000000000000000000000000000000 | | | | canconnanonnanonna | |
| | | | | | | | | | |
| | | | | | | | | | |
| | Withdrawal I | the second | | Bladder Pu | mp | | | | |
| Appeara | nce of Samp | ole: | | | | | | | |
| | Color: | *************************************** | | Clear | | | | | |
| | Turbidit | *************************************** | *********************** | Low | | | | | |
| | Sedime | ent: | | None | | | | | |
| | Other: | *************************************** | | None | | | | | |
| Laborato | ry Analysis I | Parameter: | s and Pres | ervatives: | TPH-d - 8015; | TPH-g, VOCs - | 8260; PAH | ls - 8270 | Oc sim; |
| | | | | | lead - 6020 | | | | |
| Number | and Types o | of Sample (| Containers: | 6 - 40m | nl VOAs, 2 - 1L a | amber jar, 1 - <mark>5</mark> 0 | 0ml amber | jar, 1 - 2 | 250ml HDPE |
| Sample I | Identification | Numbers: | : ES116 | [0900] | | | | | |
| Decontai | mination Pro | cedures: | Triple Rin | sed | | | | | |
| Notes: | YSI did not l | nave salini | ty paramet | er. | | | | | |
| Sampled | l by: Kirk | Markle, Je | ff Hatteme | r | | | | | |
| | Delivered to | o: <u>Ca</u> | alscience E | nvironment | | Transporters: F | edEx | | |
| Date: | 10/27/2014 | | | | | Time: 1530 | | | |
| | | | | | sing (Gallons/Lii 8"-2.61 • 10"-4. | • | | | |



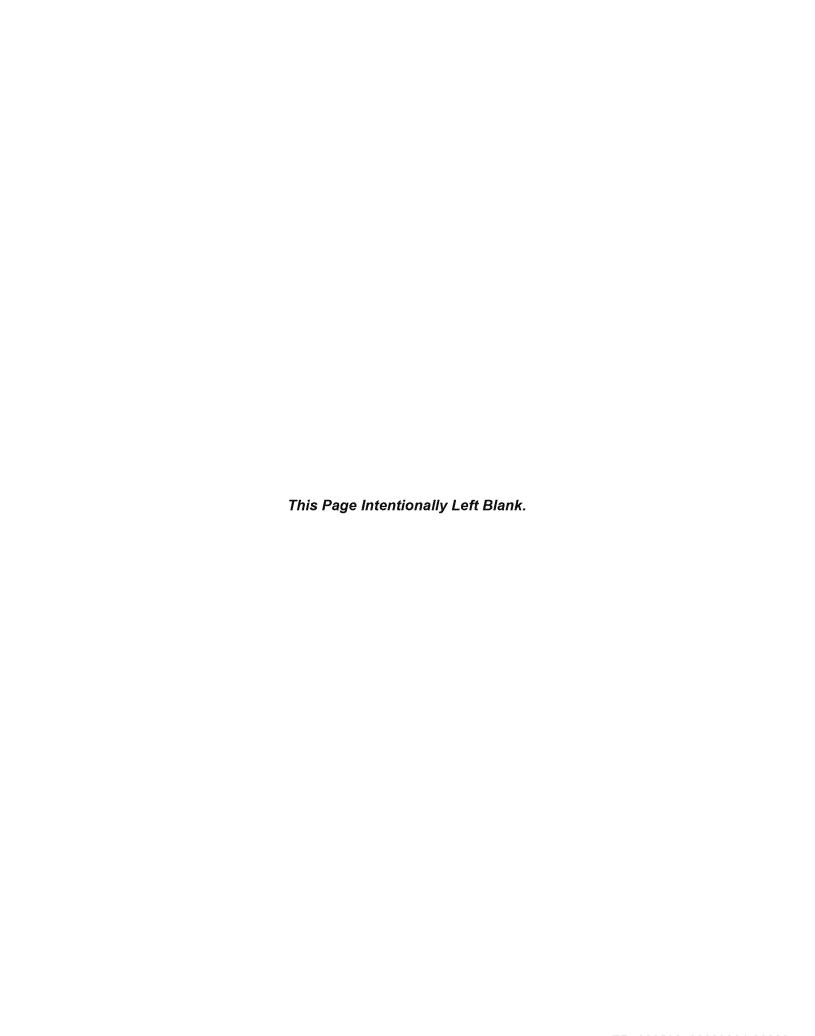


| Well ID: | RHMW05 | _ L | ocation: | Red Hil | l Bulk Fuel Stora | age Facility | Project N | o.: <u>1</u> | 12066 |
|--|--|---|------------------------|---|--|---|--|--|---|
| Initial W | ater Level: _ | 83.21 ft | | Date: | 10/28/2014 | | Time: | 1013 | |
| Total Depth of Well: Unable to Measure | | | | Personnel Involved: | | Kirk Markle, Jeff Hattemer | | | |
| Length of Saturated Zone: Unknown | | | Weath | er Conditions: | Not applicable – well is located indoors | | | ndoors | |
| Volume | of Water to b | e Removed: | 3.0 L | Method | l of Removal: | Bladder Pump | | | |
| Water L | evel After Pu | rging: | 83.21 ft | Pumpir | ng Rate: | | 0.23 L/min | | *************************************** |
| Well Pu | rge Data: | | | | | | | | |
| | Volum | ie | | nductivity | | | | | Redox (ORP) |
| Time | Remov | ed pH | <u> </u> | mS/cm) | DO (mg/l) | Temperatur | <u>e</u> Sal | <u>inity</u> | (mV) |
| 1405 | 0.0 L | 9.2 | 0 | 0.889 | 8.83 | 22.91 | | - | -211.6 |
| 1410 | 1.0 L | . 9.6 | 1 | 0.888 | 8.69 | 22.87 | | _ | -210.8 |
| 1413 | 2.0 L | 9.2 | 2 | 0.886 | 8.68 | 22.87 | | - | -210.4 |
| 1418 | 3.0 L | 9.2 | 0 | 0.886 | 8.67 | 22.85 | | - | -209.1 |
| | | | | | | | | | |
| | | | | | | | | | |
| nanananananananananananananananananana | MANAGAMAN MANAGAMANANANANANANANANANANANANANANANANA | | nanosnannan saanananan | *************************************** | *************************************** | *************************************** | 99999999999999999999999999999999999999 | ennannennannannannannannannannannannanna | *************************************** |
| | | | | | | | | | |
| | | | | | | | | | |
| | | | | | | | | | |
| • | Withdrawal N | | <u> </u> | Bladder Pu | mp | | | | |
| Appeara | ance of Samp | le: | | | | | | | |
| | Color: | *************************************** | | Clear | | | | | |
| | Turbidit | y: | | Clear | | | | | |
| | Sedime | nt: | | None | | | | | |
| | Other: | | | None | | | | | |
| Laborate | ory Analysis F | Parameters a | ind Prese | rvatives: | TPH-d - 8015 | ; TPH-g, VOCs | - 8260; PAH | ds - 8270 | Oc sim; |
| | | | | | lead - 6020 | | | | |
| Number | and Types o | f Sample Co | ntainers: | 6 - 40m | nl VOAs, 2 - 1L a | amber jar, 1 - 5 | 00ml amber | jar, 1 - 1 | IL HDPE |
| Sample | Identification | Numbers: | ES118 | [1430] | | | | | |
| • | ımination Pro | | riple Rins | ed | | | | | |
| | YSI did not h | - | | | | | | | |
| Sample | | Markle, Jeff I | | | | | | | |
| • | d Delivered to | | | vironment | al Lab | Transporters: F | edEx | | |
| Date: | 10/28/2014 | *************************************** | | | | Time: 1530 | | | ************************************** |
| | *************************************** | | | | sing (Gallons/Li | | | | |
| | | | 2"-0.16 | • 4"-0.65 • | 8"-2.61 • 10"-4. | .08 • 12"-5.87 | | | |



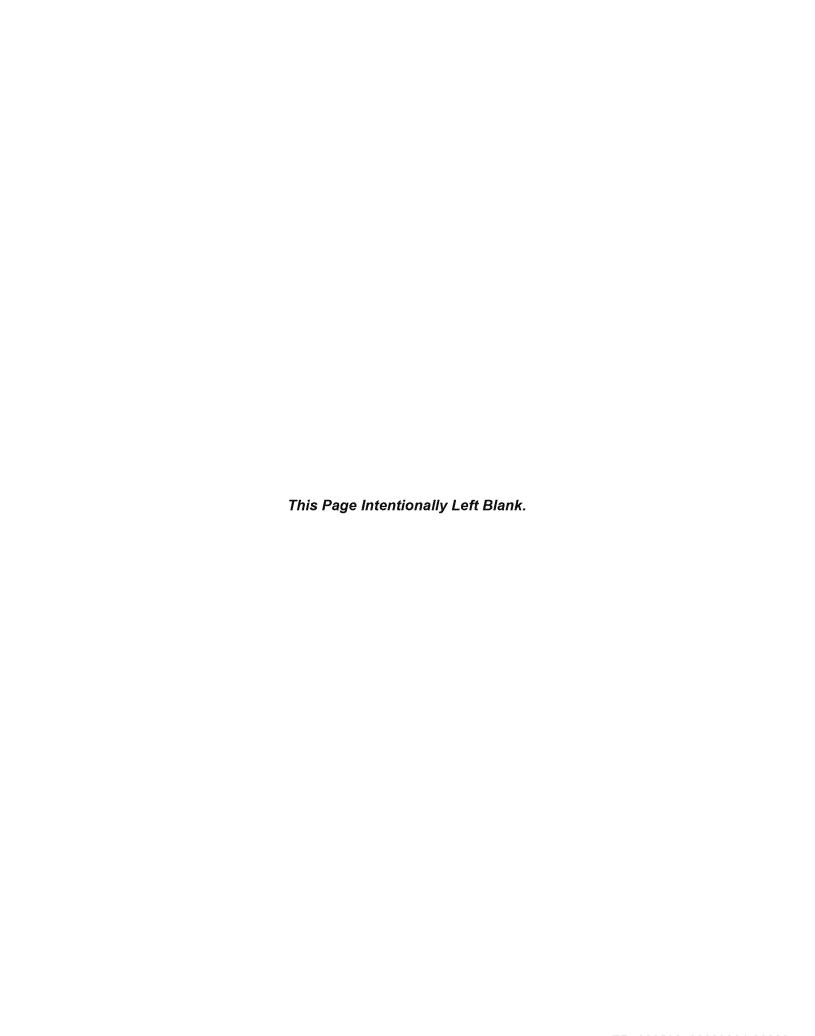


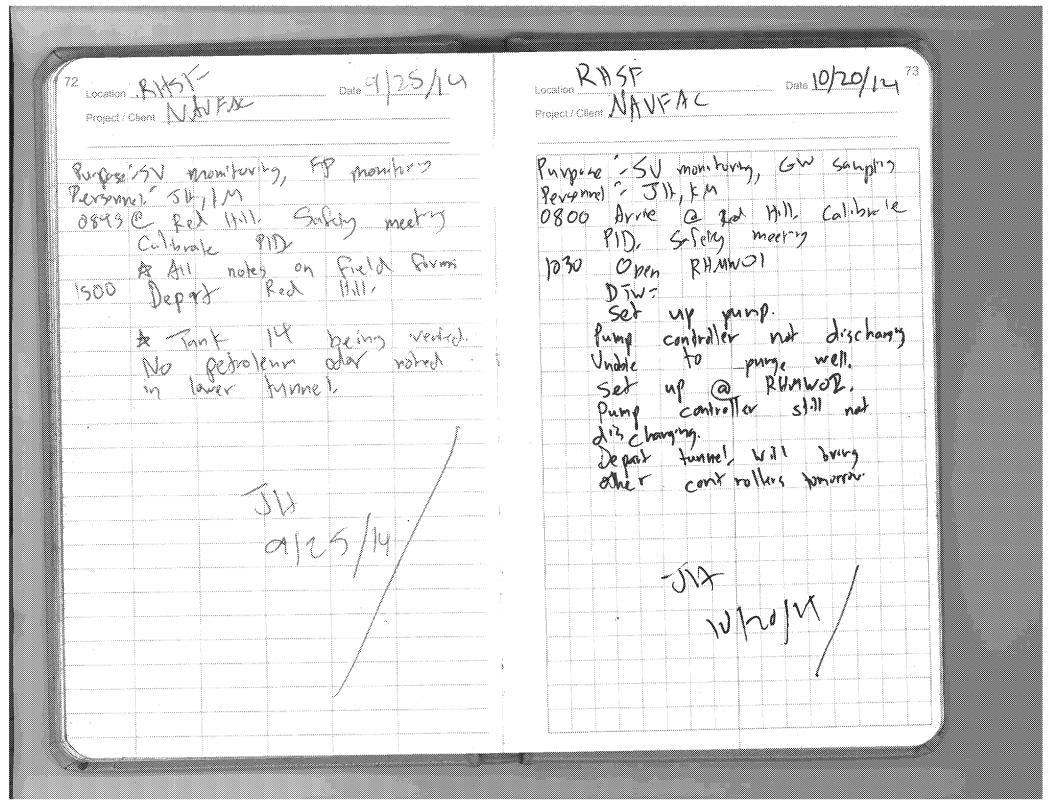
| Well ID: RHMW2254-01 | | <u>-01</u> | Location: | | Hill Bulk Fuel Stor | rage Facility | Project No.: | 112066 | | |
|--|-------------------------------------|---|------------|-------------------|--|--|---|-------------|--|--|
| Initial Wat | er Level: | 83.44 | ft | Date: _ | 10/28/2014 | | Time:11 | 10 | | |
| Total Dep | Total Depth of Well: Not applicable | | | e Per | sonnel Involved: _ | Kirk Markle, Jeff Hattemer | | | | |
| Length of Saturated Zone: Not applicable | | | ble We | ather Conditions: | Not applicab | Not applicable – well is located indoors | | | | |
| Volume of Water to be Removed: 5.0 L | | | | | hod of Removal: _ | Е | Bladder Pump | | | |
| Water Level After Purging: 83.44 ft | | | | | nping Rate: | 0.50 L/min | | | | |
| Well Purg | | | | | | | | | | |
| | Volum | | | Conductiv | • | | - u u | Redox (ORP) | | |
| Time | _ Remov | <u>ed</u> | pH | (mS/cm | <u> </u> | Temperatur | e Salinity | (mV) | | |
| 1117 | 0.0 L | | 9.60 | 0.613 | 13.00 | 21.97 | _ | 201.7 | | |
| 1119 | 1.0 L | *************************************** | 9.55 | 0.606 | 10.29 | 21.59 | | 200.3 | | |
| 1121 | 2.0 L | | 9.65 | 0.603 | 9.57 | 21.45 | <u> </u> | 197.5 | | |
| 1123 | 3.0 L | | 9.47 | 0.603 | 9.43 | 21.38 | | 196.2 | | |
| 1125 | 4.0 L | | 9.44 | 0.603 | 9.23 | 21.28 | | 195.0 | | |
| 1127 | 5.0 L | | 9.43 | 0.603 | 9.17 | 21.28 | | | | |
| | | | | | | | | | | |
| Sample W | /ithdrawal M | lethod: | | Bladder | Pump | _ | | | | |
| Appearan | ce of Samp | e: | | | | | | | | |
| | Color: | | | Clear | | _ | | | | |
| | Turbidity | / : | | Clear | | | | | | |
| | Sedime | nt: | | None | | _ | | | | |
| | Other: | *************************************** | | None | | | | | | |
| Laborator | y Analysis P | aramete | ers and Pi | reservatives | s: <u>TPH-d - 8015</u> lead - 200.8 | 5; TPH-g, VOCs | - 8260; PAHs - 8 | 3270c sim; | | |
| Number a | nd Types of | Sample | Containe | ers: 6-4 | 40ml VOAs, 2 - 1L | amber jar, 1 - 5 | 00ml amber jar, 1 | 1 - 1L HDPE | | |
| Sample Id | lentification | Number | s: ES | 117 [1200] | | | | | | |
| Decontarr | nination Prod | cedures | | | | | | | | |
| | 'SI did not h | | | | | | *************************************** | | | |
| Sampled | | | Jeff Hatte | | | | | | | |
| • | Delivered to | | | e Environm | ental Lab | Transporters: F | edEx | | | |
| Date: 1 | 0/28/2014 | | | | NATIONAL PROPERTY OF THE PROPE | Time: 1530 | *************************************** | | | |
| | | | | | Casing (Gallons/L 55 • 8"-2.61 • 10"-4 | • | | | | |

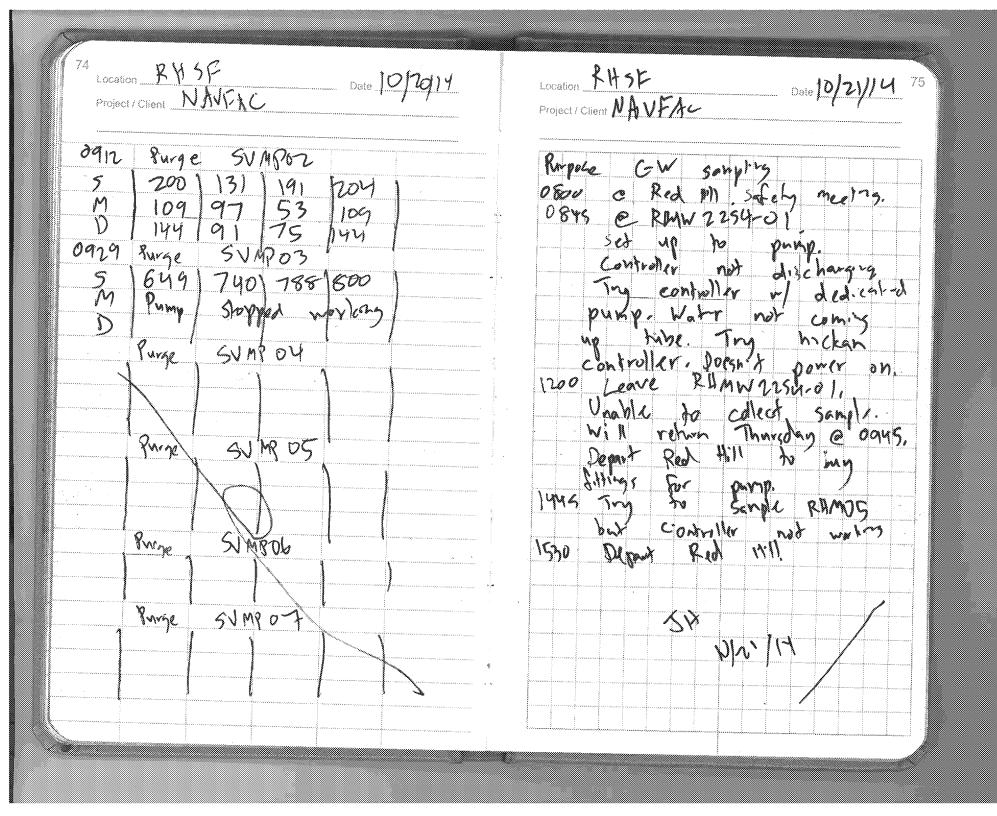


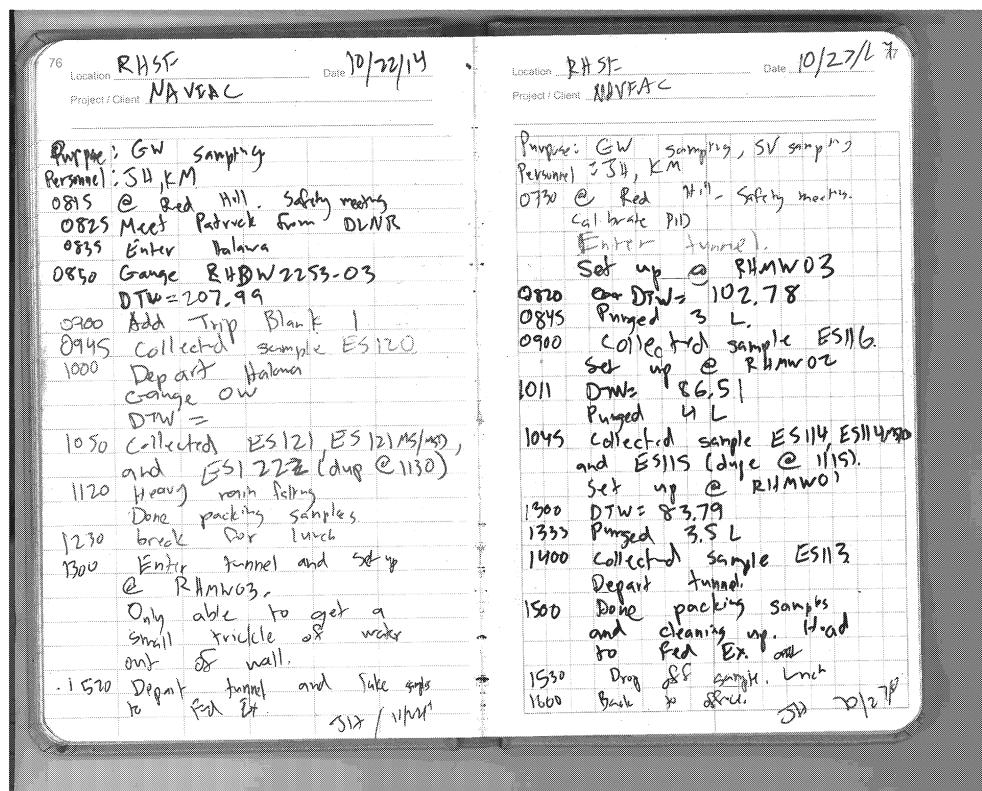
APPENDIX B

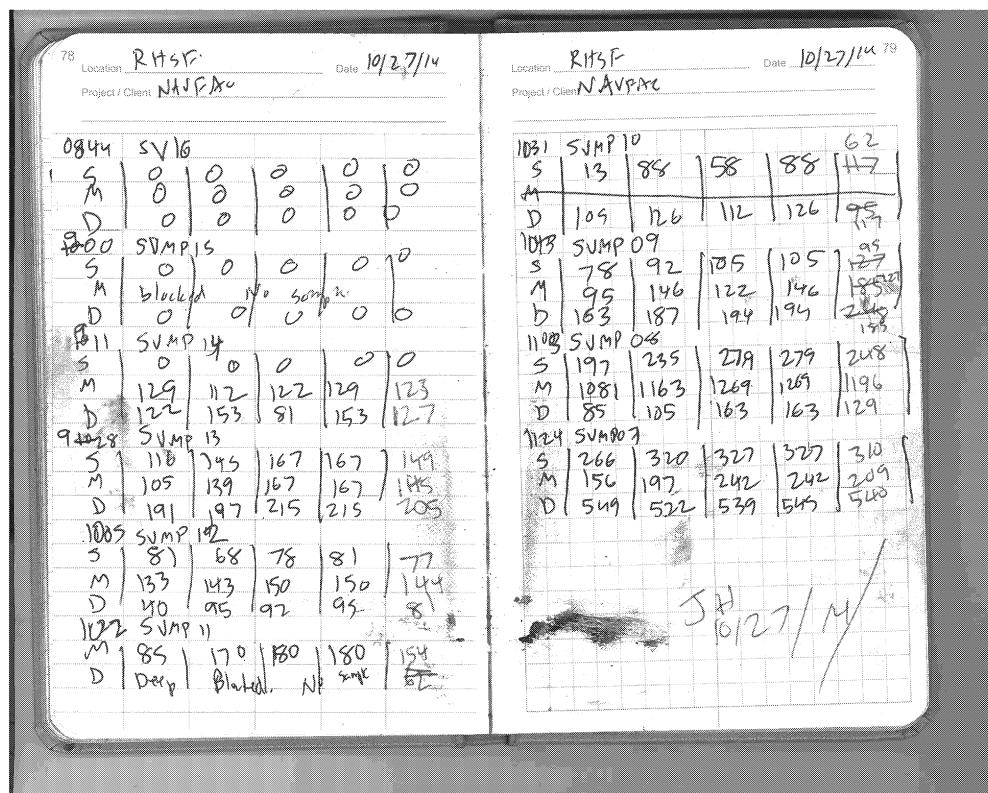
Field Notes

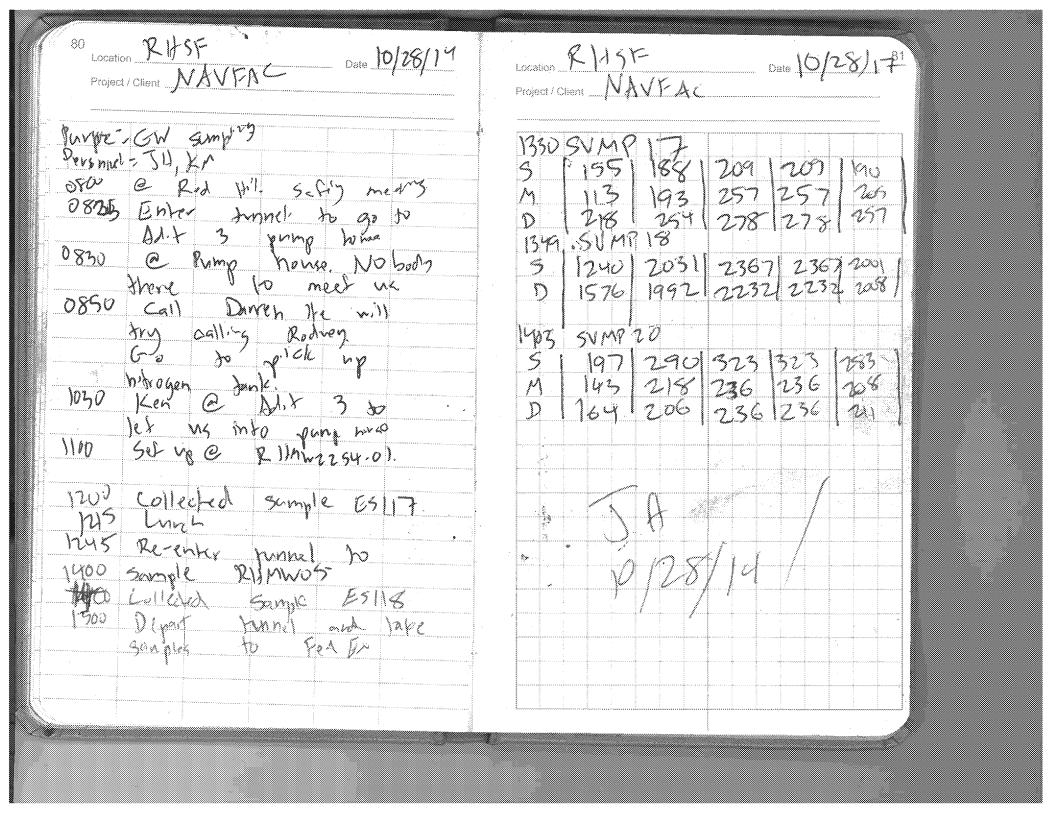


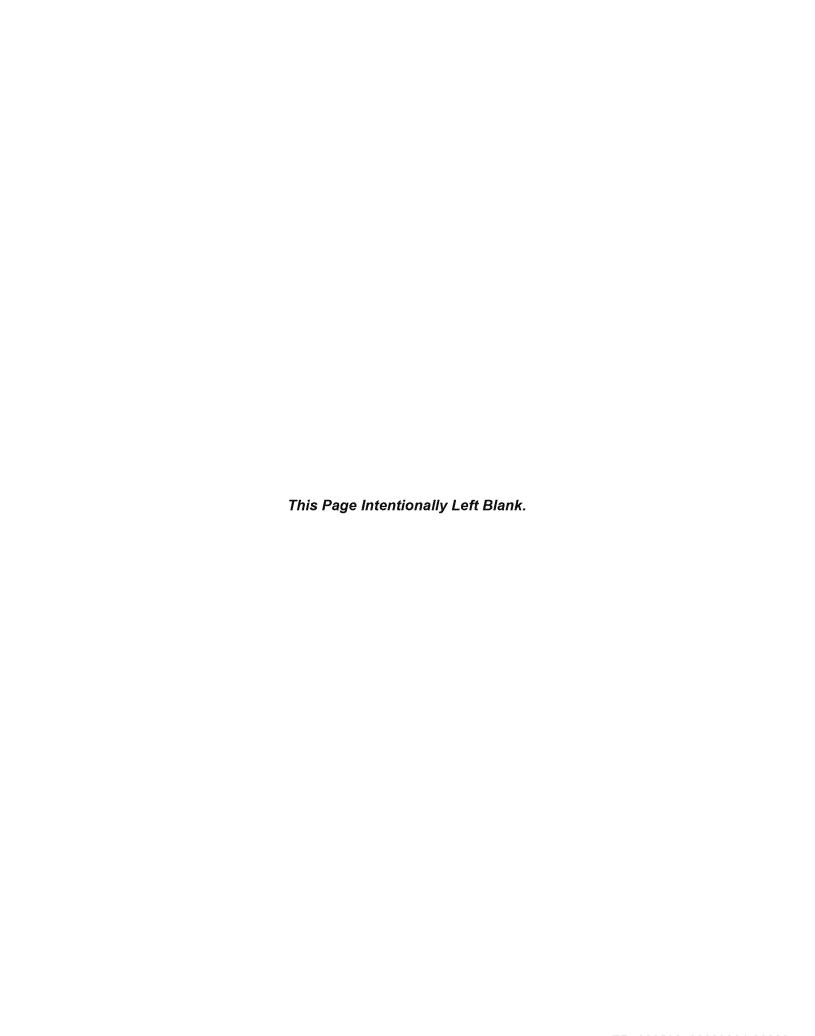




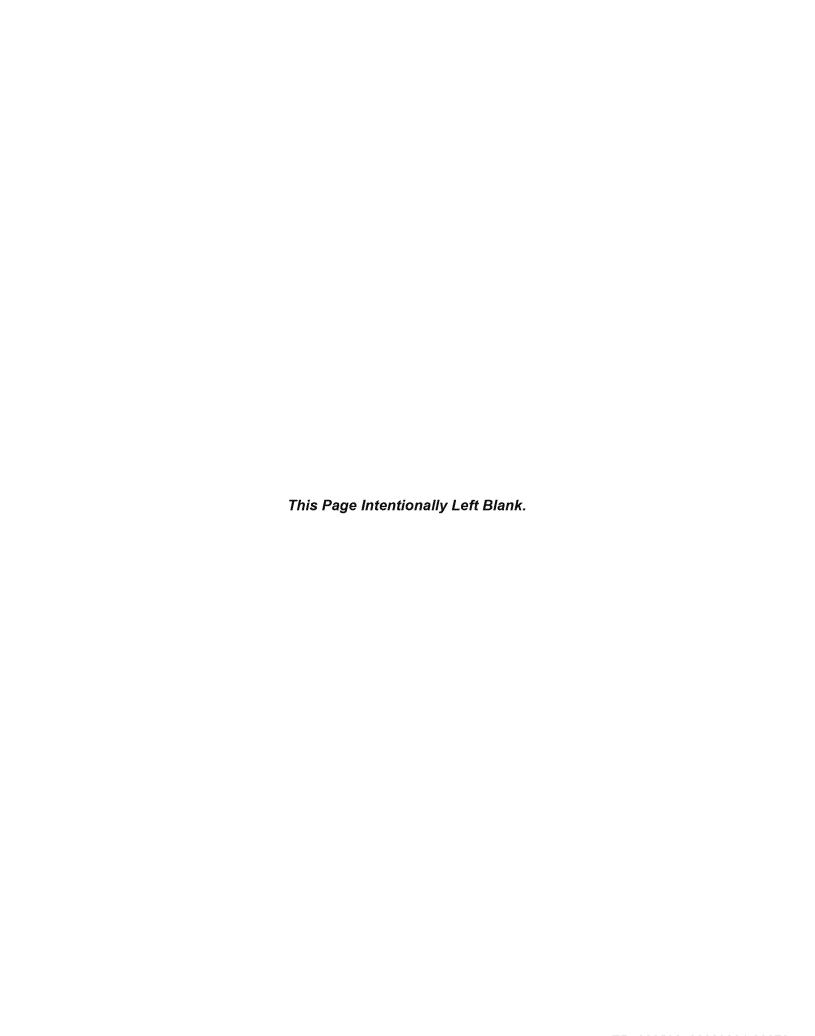








APPENDIX C Laboratory Reports





Calscience



WORK ORDER NUMBER: 14-10-2131

The difference is service



AIR SOIL WATER | MARINE CHEMISTRY

Analytical Report For

Client: Environmental Science International, Inc.

Client Project Name: Red Hill LTM 112066

Attention: Jeff Hattemer

354 Uluniu Street, Suite 304 Kailua, HI 96734-2500

Rand Ollo

Approved for release on 11/04/2014 by: Richard Villafania Project Manager



Email your FM F

Resulting

Eurofins Calscience, Inc. (Calscience) certifies that the test results provided in this report meet all NELAC requirements for parameters for which accreditation is required or available. Any exceptions to NELAC requirements are noted in the case narrative. The original report of subcontracted analyses, if any, is attached to this report. The results in this report are limited to the sample(s) tested and any reproduction thereof must be made in its entirety. The client or recipient of this report is specifically prohibited from making material changes to said report and, to the extent that such changes are made, Calscience is not responsible, legally or otherwise. The client or recipient agrees to indemnify Calscience for any defense to any litigation which may arise.

744) Enralo Wey Gurren Scree (A.9.264) (4.2.) • TEL 7131 295 (294 • F2.677) 4 894 7591 • gross decimandos

NELAP DI COURCA : ACLASS DODELAP ID AGE 1864 (SOULCITIOS 2005). CSD ACLD 10:00 (I SCACIAL DI RICALITA)



Contents

| Client Project Name: | Red Hill LTM 112066 |
|----------------------|---------------------|
| Work Order Number: | 14-10-2131 |

| 1 | Work Order Narrative | 3 |
|---|--|------------------------|
| 2 | Client Sample Data. 2.1 EPA 8015B (M) TPH Diesel (Aqueous). 2.2 EPA 6020 ICP/MS Metals (Aqueous). 2.3 EPA 8270C SIM PAHs (Aqueous). 2.4 GC/MS GRO/EPA 8260B Volatile Organics (Aqueous). | 4 4 6 7 12 |
| 3 | Quality Control Sample Data. 3.1 MS/MSD. 3.2 PDS/PDSD. 3.3 LCS/LCSD. | 24 24 29 30 |
| 4 | Sample Analysis Summary | 35 |
| 5 | Glossary of Terms and Qualifiers | 36 |
| 6 | Chain-of-Custody/Sample Receipt Form | 37 |



Work Order Narrative

Work Order: 14-10-2131 Page 1 of 1

Condition Upon Receipt:

Samples were received under Chain-of-Custody (COC) on 10/28/14. They were assigned to Work Order 14-10-2131.

Unless otherwise noted on the Sample Receiving forms all samples were received in good condition and within the recommended EPA temperature criteria for the methods noted on the COC. The COC and Sample Receiving Documents are integral elements of the analytical report and are presented at the back of the report.

Holding Times:

All samples were analyzed within prescribed holding times (HT) and/or in accordance with the Calscience Sample Acceptance Policy unless otherwise noted in the analytical report and/or comprehensive case narrative, if required.

Any parameter identified in 40CFR Part 136.3 Table II that is designated as "analyze immediately" with a holding time of <= 15 minutes (40CFR-136.3 Table II, footnote 4), is considered a "field" test and the reported results will be qualified as being received outside of the stated holding time unless received at the laboratory within 15 minutes of the collection time.

Quality Control:

All quality control parameters (QC) were within established control limits except where noted in the QC summary forms or described further within this report.

Additional Comments:

Air - Sorbent-extracted air methods (EPA TO-4A, EPA TO-10, EPA TO-13A, EPA TO-17): Analytical results are converted from mass/sample basis to mass/volume basis using client-supplied air volumes.

New York NELAP air certification does not certify for all reported methods and analytes, reference the accredited items here: http://www.calscience.com/PDF/New_York.pdf

Solid - Unless otherwise indicated, solid sample data is reported on a wet weight basis, not corrected for % moisture. All QC results are always reported on a wet weight basis.

Subcontractor Information:

Unless otherwise noted below (or on the subcontract form), no samples were subcontracted.



10/28/14 Date Received: Environmental Science International, Inc. Work Order: 14-10-2131 354 Uluniu Street, Suite 304 Preparation: **EPA 3510C** Kailua, HI 96734-2500 Method: EPA 8015B (M) Units: ug/L Project: Red Hill LTM 112066 Page 1 of 2 Client Sample Number Lab Sample Date/Time Matrix Instrument Date Date/Time QC Batch ID Number Collected Prepared Analyzed ES113 14-10-2131-2-H 10/27/14 Aqueous GC 45 10/30/14 10/31/14 141030B06 Comment(s): - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag. - TPH as Diesel is quantified in the carbon range C10-C28. Parameter Result DL LOD LOQ DF Qualifiers TPH as Diesel 120 11 12 25 1.00 HD Surrogate Rec. (%) **Control Limits** Qualifiers n-Octacosane 51-141 91 10/27/14 10:45 10/31/14 09:59 ES114 10/30/14 141030B06 14-10-2131-3-H Aqueous GC 45 - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag. Comment(s): - TPH as Diesel is quantified in the carbon range C10-C28. <u>LOD</u> Parameter Result LOQ <u>DF</u> Qualifiers TPH as Diesel 2000 11 12 25 1.00 HD Qualifiers Surrogate Rec. (%) Control Limits 51-141 n-Octacosane 92 ES115 14-10-2131-4-H 10/27/14 GC 45 10/30/14 10/31/14 141030B06 Aqueous 10:41 Comment(s): - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag. - TPH as Diesel is quantified in the carbon range C10-C28. **Parameter** Result DL LOD LOQ DF Qualifiers 2000 TPH as Diesel 11 12 25 1.00 HDSurrogate Rec. (%) **Control Limits** Qualifiers 51-141 n-Octacosane 91 ES116 14-10-2131-5-H 10/27/14 GC 45 10/30/14 141030B06 Aqueous 10/31/14 09:00 10:58 Comment(s): - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag. - TPH as Diesel is quantified in the carbon range C10-C28. LOQ <u>DF</u> Parameter Result DL LOD Qualifiers TPH as Diesel 12 25 1.00 HD **Control Limits** Qualifiers Surrogate Rec. (%)

51-141

65

n-Octacosane



Environmental Science International, Inc.

Date Received: 10/28/14

Work Order: 14-10-2131

Kailua, HI 96734-2500 Preparation: EPA 3510C

Method: EPA 8015B (M)

Units: ug/L

Project: Red Hill LTM 112066 Page 2 of 2

| Client Sample Number | Lab Sample Number | Date/Time MacCollected | atrix Instrume | nt Date Prepared | Date/Time Analyzed | QC Batch ID |
|---|----------------------|------------------------|-------------------|---------------------|-----------------------|-------------------|
| Method Blank | 099-15-516-206 | N/A A | queous GC 45 | 10/30/14 | 10/31/14 07:03 | 141030B06 |
| Comment(s): - Results were evaluated to | the MDL (DL), cond | entrations >= to the | MDL (DL) but < RL | (LOQ), if found, a | re qualified with | a "J" flag. |
| <u>Parameter</u> | <u>Result</u> | <u>DL</u> | LOD | LOQ | <u>DF</u> | <u>Qualifiers</u> |
| TPH as Diesel | <12 | 11 | 12 | 25 | 1.00 | U |
| | | | | | | |
| Surrogate | Rec. (%) | Control Limits | <u>Qualifiers</u> | | | |
| n-Octacosane | 117 | 51-141 | | | | |





10/28/14 Date Received: Environmental Science International, Inc. Work Order: 14-10-2131 354 Uluniu Street, Suite 304 Preparation: EPA 3005A Filt. Kailua, HI 96734-2500 Method: EPA 6020 Units: ug/L Project: Red Hill LTM 112066 Page 1 of 1 Client Sample Number Lab Sample Date/Time Matrix Instrument Date Date/Time QC Batch ID Number Collected Prepared Analyzed ES113 14-10-2131-2-G 10/27/14 Aqueous ICP/MS 04 10/29/14 10/30/14 141029L03D - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag. Comment(s): <u>Parameter</u> Result DL **LOD** LOQ DF Qualifiers Lead 0.0976 0.0898 0.200 1.00 1.00 ES114 10/27/14 10/30/14 14-10-2131-3-G Aqueous ICP/MS 04 10/29/14 141029L03D 10:45 23:32 Comment(s): - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag. LOD LOQ Parameter Result <u>DL</u> DF Qualifiers U Lead < 0.200 0.0898 0.200 1.00 1.00 14-10-2131-4-G 10/27/14 11:15 ES115 ICP/MS 04 10/29/14 10/30/14 141029L03D Aqueous 23:39 - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag. Comment(s): <u>Parameter</u> Result <u>DL</u> <u>LOD</u> LOQ <u>DF</u> Qualifiers 0.165 0.0898 0.200 1.00 Lead 1.00 ES116 14-10-2131-5-G 10/27/14 ICP/MS 04 10/29/14 10/30/14 141029L03D Aqueous 09:00 23:43 Comment(s): - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag. <u>Parameter</u> Result <u>DL</u> <u>LOD</u> LOQ <u>DF</u> Qualifiers Lead < 0.200 0.0898 0.200 1.00 1.00 U

| Method Blank | 099-14-497-1 | 01 N/A | Aqueous II | CP/MS 04 10/29 | /14 10/30/1 22:58 | 4 141029L03D |
|------------------|---|------------------|---------------------|------------------------|----------------------|-------------------|
| Comment(s): | - Results were evaluated to the MDL (DL), | concentrations > | = to the MDL (DL) b | out < RL (LOQ), if for | ınd, are qualified v | with a "J" flag. |
| <u>Parameter</u> | Result | <u>DL</u> | LOD | LOQ | <u>DF</u> | <u>Qualifiers</u> |
| Lead | <0.200 | 0.0898 | 3 0.200 | 1.00 | 1.00 | U |



Environmental Science International, Inc.

354 Uluniu Street, Suite 304 Kailua, HI 96734-2500 Date Received: Work Order: Preparation:

EPA 3510C EPA 8270C SIM PAHs

Method: Units:

ug/L

10/28/14

14-10-2131

Project: Red Hill LTM 112066

Page 1 of 5

| Client Sample Number | Lab Sample Number | Date/Time N | Matrix | Instrument | | Date/Time QC Batch ID Analyzed |
|--------------------------------------|----------------------|----------------------|-------------|---------------|---------------------|-----------------------------------|
| ES113 | 14-10-2131-2-1 | 10/27/14 A 14:00 | Aqueous | GC/MS AAA | | 10/31/14 141030L08 20:19 |
| Comment(s): - Results were evaluated | to the MDL (DL), con | centrations >= to th | e MDL (DL) | but < RL (LOQ |), if found, are qu | ualified with a "J" flag. |
| <u>Parameter</u> | Result | <u>DL</u> | LOD | LOQ | <u>DF</u> | <u>Qualifiers</u> |
| Naphthalene | <0.052 | 0.035 | 0.052 | 0.21 | 1.0 | 0 U |
| 2-Methylnaphthalene | <0.052 | 0.048 | 0.052 | 0.21 | 1.0 | 0 U |
| 1-Methylnaphthalene | <0.10 | 0.054 | 0.10 | 0.21 | 1.0 | 0 U |
| Acenaphthylene | <0.052 | 0.046 | 0.052 | 0.21 | 1.0 | 0 U |
| Acenaphthene | <0.052 | 0.028 | 0.052 | 0.21 | 1.0 | 0 U |
| Fluorene | <0.052 | 0.044 | 0.052 | 0.21 | 1.0 | 0 U |
| Phenanthrene | <0.052 | 0.028 | 0.052 | 0.21 | 1.0 | 0 U |
| Anthracene | <0.052 | 0.030 | 0.052 | 0.21 | 1.0 | 0 U |
| Fluoranthene | <0.052 | 0.049 | 0.052 | 0.21 | 1.0 | 0 U |
| Pyrene | <0.052 | 0.021 | 0.052 | 0.21 | 1.0 | 0 U |
| Benzo (a) Anthracene | <0.052 | 0.034 | 0.052 | 0.21 | 1.0 | 0 U |
| Chrysene | <0.052 | 0.026 | 0.052 | 0.21 | 1.0 | 0 U |
| Benzo (k) Fluoranthene | <0.052 | 0.032 | 0.052 | 0.21 | 1.0 | 0 U |
| Benzo (b) Fluoranthene | <0.052 | 0.018 | 0.052 | 0.21 | 1.0 | 0 U |
| Benzo (a) Pyrene | <0.052 | 0.023 | 0.052 | 0.21 | 1.0 | 0 U |
| Indeno (1,2,3-c,d) Pyrene | <0.052 | 0.022 | 0.052 | 0.21 | 1.0 | 0 U |
| Dibenz (a,h) Anthracene | <0.052 | 0.049 | 0.052 | 0.21 | 1.0 | 0 U |
| Benzo (g,h,i) Perylene | <0.10 | 0.085 | 0.10 | 0.21 | 1.0 | 0 U |
| Surrogate | Rec. (%) | Control Limit | ts Qualifie | <u>ers</u> | | |
| Nitrobenzene-d5 | 49 | 28-139 | | | | |
| 2-Fluorobiphenyl | 65 | 33-144 | | | | |
| p-Terphenyl-d14 | 70 | 23-160 | | | | |



p-Terphenyl-d14

Analytical Report

Environmental Science International, Inc.

Date Received: 10/28/14

354 Uluniu Street, Suite 304

Work Order: 14-10-2131

Kailua, HI 96734-2500 Preparation: EPA 3510C Method: EPA 8270C SIM PAHs

Units: ug/L

Project: Red Hill LTM 112066 Page 2 of 5

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|--|-----------------------|------------------------|--------------------|-----------------|------------------|-----------------------|-------------------|
| ES114 | 14-10-2131-3-J | 10/27/14 10:45 | Aqueous | GC/MS AAA | 10/30/14 | 10/31/14 19:10 | 141030L08 |
| Comment(s): - Results were evaluated t | to the MDL (DL), cond | centrations >= to t | he MDL (DL |) but < RL (LOC | (a), if found, a | are qualified with | a "J" flag. |
| <u>Parameter</u> | <u>Result</u> | <u>DL</u> | <u>LOD</u> | LOC | Ī | <u>DF</u> | <u>Qualifiers</u> |
| Acenaphthylene | <0.047 | 0.042 | 0.047 | 0.19 | 1 | 1.00 | U |
| Acenaphthene | 0.53 | 0.025 | 0.047 | 0.19 | 1 | 1.00 | |
| Fluorene | <0.047 | 0.040 | 0.047 | 0.19 | 1 | 1.00 | U |
| Phenanthrene | <0.047 | 0.026 | 0.047 | 0.19 | 1 | 1.00 | U |
| Anthracene | <0.047 | 0.027 | 0.047 | 0.19 | 1 | 1.00 | U |
| Fluoranthene | <0.047 | 0.044 | 0.047 | 0.19 | 1 | 1.00 | U |
| Pyrene | <0.047 | 0.019 | 0.047 | 0.19 | 1 | 1.00 | U |
| Benzo (a) Anthracene | <0.047 | 0.031 | 0.047 | 0.19 | 1 | 1.00 | U |
| Chrysene | <0.047 | 0.024 | 0.047 | 0.19 | 1 | 1.00 | U |
| Benzo (k) Fluoranthene | <0.047 | 0.029 | 0.047 | 0.19 | 1 | 1.00 | U |
| Benzo (b) Fluoranthene | <0.047 | 0.017 | 0.047 | 0.19 | 1 | 1.00 | U |
| Benzo (a) Pyrene | <0.047 | 0.021 | 0.047 | 0.19 | 1 | 1.00 | U |
| Indeno (1,2,3-c,d) Pyrene | <0.047 | 0.020 | 0.047 | 0.19 | 1 | 1.00 | U |
| Dibenz (a,h) Anthracene | <0.047 | 0.045 | 0.047 | 0.19 | 1 | 1.00 | U |
| Benzo (g,h,i) Perylene | <0.095 | 0.078 | 0.095 | 0.19 | 1 | 1.00 | U |
| | | | | | | | |
| Surrogate | Rec. (%) | Control Limi | <u>its Qualifi</u> | iers | | | |
| Nitrobenzene-d5 | 66 | 28-139 | | | | | |
| 2-Fluorobiphenyl | 65 | 33-144 | | | | | |

| ES114 | 14-10-2131-3-J | 10/27/14 A 10:45 | kqueous GC | /MS AAA 10/30/1 | 4 11/03/14 18:28 | 141030L08 |
|---|---------------------|----------------------|-------------------|-------------------|-----------------------|-------------------|
| Comment(s): - Results were evaluated to | the MDL (DL), conce | entrations >= to the | e MDL (DL) but | RL (LOQ), if foun | d, are qualified with | a "J" flag. |
| <u>Parameter</u> | <u>Result</u> | <u>DL</u> | LOD | <u>LOQ</u> | <u>DF</u> | <u>Qualifiers</u> |
| Naphthalene | 140 | 0.64 | 0.95 | 3.8 | 20.0 | |
| 2-Methylnaphthalene | 43 | 0.88 | 0.95 | 3.8 | 20.0 | |
| 1-Methylnaphthalene | 59 | 0.98 | 1.9 | 3.8 | 20.0 | |
| Surrogate | Rec. (%) | Control Limits | <u>Qualifiers</u> | | | |
| Nitrobenzene-d5 | 60 | 28-139 | | | | |
| 2-Fluorobiphenyl | 80 | 33-144 | | | | |
| p-Terphenyl-d14 | 83 | 23-160 | | | | |

23-160

68

10/28/14



p-Terphenyl-d14

Analytical Report

Date Received: Environmental Science International, Inc. 354 Uluniu Street, Suite 304 Work Order: 14-10-2131

Preparation: EPA 3510C Kailua, HI 96734-2500 Method: EPA 8270C SIM PAHs

> Units: ug/L

Project: Red Hill LTM 112066 Page 3 of 5

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|--|----------------------|------------------------|--------------------|-----------------|------------------|-----------------------|-------------------|
| ES115 | 14-10-2131-4-1 | 10/27/14 11:15 | Aqueous | GC/MS AAA | 10/30/14 | 10/31/14 19:33 | 141030L08 |
| Comment(s): - Results were evaluated t | o the MDL (DL), con | centrations >= to the | he MDL (DL |) but < RL (LOC | (a), if found, a | are qualified with | a "J" flag. |
| <u>Parameter</u> | <u>Result</u> | <u>DL</u> | <u>LOD</u> | LOC | Ī | <u>DF</u> | <u>Qualifiers</u> |
| Acenaphthylene | <0.047 | 0.042 | 0.047 | 0.19 | 1 | 1.00 | U |
| Acenaphthene | 0.53 | 0.025 | 0.047 | 0.19 | 1 | 1.00 | |
| Fluorene | <0.047 | 0.040 | 0.047 | 0.19 | 1 | 1.00 | U |
| Phenanthrene | <0.047 | 0.026 | 0.047 | 0.19 | 1 | 1.00 | U |
| Anthracene | <0.047 | 0.027 | 0.047 | 0.19 | 1 | 1.00 | U |
| Fluoranthene | <0.047 | 0.044 | 0.047 | 0.19 | 1 | 1.00 | U |
| Pyrene | <0.047 | 0.019 | 0.047 | 0.19 | 1 | 1.00 | U |
| Benzo (a) Anthracene | <0.047 | 0.031 | 0.047 | 0.19 | 1 | 1.00 | U |
| Chrysene | <0.047 | 0.023 | 0.047 | 0.19 | 1 | 1.00 | U |
| Benzo (k) Fluoranthene | <0.047 | 0.029 | 0.047 | 0.19 | 1 | 1.00 | U |
| Benzo (b) Fluoranthene | <0.047 | 0.017 | 0.047 | 0.19 | 1 | 1.00 | U |
| Benzo (a) Pyrene | <0.047 | 0.021 | 0.047 | 0.19 | 1 | 1.00 | U |
| Indeno (1,2,3-c,d) Pyrene | <0.047 | 0.020 | 0.047 | 0.19 | 1 | 1.00 | U |
| Dibenz (a,h) Anthracene | <0.047 | 0.045 | 0.047 | 0.19 | 1 | 1.00 | U |
| Benzo (g,h,i) Perylene | <0.095 | 0.077 | 0.095 | 0.19 |) | 1.00 | U |
| | | | | | | | |
| Surrogate | Rec. (%) | Control Limi | <u>its Qualifi</u> | <u>iers</u> | | | |
| Nitrobenzene-d5 | 70 | 28-139 | | | | | |
| 2-Fluorobiphenyl | 69 | 33-144 | | | | | |

| ES115 | 14-10-2131-4-1 | 10/27/14 A 11:15 | queous GC | MS AAA 10/30/14 | 11/03/14 18:52 | 141030L08 |
|---|--------------------|----------------------|----------------|----------------------|-----------------------|-------------------|
| Comment(s): - Results were evaluated to | the MDL (DL), conc | entrations >= to the | e MDL (DL) but | < RL (LOQ), if found | d, are qualified with | a "J" flag. |
| <u>Parameter</u> | <u>Result</u> | <u>DL</u> | LOD | LOQ | <u>DF</u> | <u>Qualifiers</u> |
| Naphthalene | 130 | 0.64 | 0.95 | 3.8 | 20.0 | |
| 2-Methylnaphthalene | 36 | 0.88 | 0.95 | 3.8 | 20.0 | |
| 1-Methylnaphthalene | 54 | 0.98 | 1.9 | 3.8 | 20.0 | |
| Surrogate | Rec. (%) | Control Limits | Qualifiers | | | |
| Nitrobenzene-d5 | 53 | 28-139 | | | | |
| 2-Fluorobiphenyl | 78 | 33-144 | | | | |
| p-Terphenyl-d14 | 80 | 23-160 | | | | |

23-160

73

10/28/14

14-10-2131

EPA 3510C



Analytical Report

Environmental Science International, Inc.

Date Received:

Work Order:

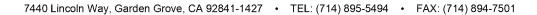
Kailua, HI 96734-2500 Preparation:

Method: EPA 8270C SIM PAHs

Units: ug/L

Project: Red Hill LTM 112066 Page 4 of 5

| Client Sample Number | Lab Sample Number | Date/Time I Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|--|----------------------|--------------------------|-------------------|-----------------|------------------|-----------------------|-------------------|
| ES116 | 14-10-2131-5-1 | 10/27/14 / 09:00 | Aqueous | GC/MS AAA | 10/30/14 | 10/31/14 19:56 | 141030L08 |
| Comment(s): - Results were evaluated t | o the MDL (DL), con | centrations >= to th | ne MDL (DL |) but < RL (LOC |), if found, a | re qualified with a | a "J" flag. |
| <u>Parameter</u> | <u>Result</u> | <u>DL</u> | <u>LOD</u> | LOC | <u>)</u> | <u>DF</u> | <u>Qualifiers</u> |
| Naphthalene | <0.054 | 0.036 | 0.054 | 0.21 | | 1.00 | U |
| 2-Methylnaphthalene | <0.054 | 0.050 | 0.054 | 0.21 | | 1.00 | U |
| 1-Methylnaphthalene | <0.11 | 0.055 | 0.11 | 0.21 | | 1.00 | U |
| Acenaphthylene | <0.054 | 0.048 | 0.054 | 0.21 | | 1.00 | U |
| Acenaphthene | <0.054 | 0.029 | 0.054 | 0.21 | | 1.00 | U |
| Fluorene | <0.054 | 0.046 | 0.054 | 0.21 | | 1.00 | U |
| Phenanthrene | <0.054 | 0.029 | 0.054 | 0.21 | | 1.00 | U |
| Anthracene | <0.054 | 0.031 | 0.054 | 0.21 | | 1.00 | U |
| Fluoranthene | <0.054 | 0.050 | 0.054 | 0.21 | | 1.00 | U |
| Pyrene | <0.054 | 0.022 | 0.054 | 0.21 | | 1.00 | U |
| Benzo (a) Anthracene | <0.054 | 0.035 | 0.054 | 0.21 | | 1.00 | U |
| Chrysene | <0.054 | 0.027 | 0.054 | 0.21 | | 1.00 | U |
| Benzo (k) Fluoranthene | <0.054 | 0.033 | 0.054 | 0.21 | | 1.00 | U |
| Benzo (b) Fluoranthene | <0.054 | 0.019 | 0.054 | 0.21 | | 1.00 | U |
| Benzo (a) Pyrene | <0.054 | 0.024 | 0.054 | 0.21 | | 1.00 | U |
| Indeno (1,2,3-c,d) Pyrene | <0.054 | 0.023 | 0.054 | 0.21 | | 1.00 | U |
| Dibenz (a,h) Anthracene | <0.054 | 0.051 | 0.054 | 0.21 | | 1.00 | U |
| Benzo (g,h,i) Perylene | <0.11 | 0.088 | 0.11 | 0.21 | | 1.00 | U |
| Surrogate | Rec. (%) | Control Limit | <u>ts Qualifi</u> | iers | | | |
| Nitrobenzene-d5 | 59 | 28-139 | | | | | |
| 2-Fluorobiphenyl | 67 | 33-144 | | | | | |
| p-Terphenyl-d14 | 74 | 23-160 | | | | | |





Environmental Science International, Inc.

354 Uluniu Street, Suite 304 Kailua, HI 96734-2500 Date Received:

10/28/14 14-10-2131

Work Order: Preparation:

EPA 3510C

Method:

EPA 8270C SIM PAHs

Units:

ug/L

Project: Red Hill LTM 112066

Page 5 of 5

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|--------------------------------------|----------------------|------------------------|--------------------|-----------------|------------------|-----------------------|-------------------|
| Method Blank | 099-15-148-67 | N/A | Aqueous | GC/MS AAA | 10/30/14 | 10/31/14 17:17 | 141030L08 |
| Comment(s): - Results were evaluated | to the MDL (DL), cor | ncentrations >= to th | he MDL (DL |) but < RL (LOC |), if found, are | qualified with a | "J" flag. |
| <u>Parameter</u> | <u>Result</u> | <u>DL</u> | <u>LOD</u> | LOC | <u>]</u> | <u>DF</u> | <u>Qualifiers</u> |
| Naphthalene | <0.050 | 0.034 | 0.050 | 0.20 | 1 | 1.00 | U |
| 2-Methylnaphthalene | <0.050 | 0.046 | 0.050 | 0.20 | 1 | 1.00 | U |
| 1-Methylnaphthalene | <0.10 | 0.052 | 0.10 | 0.20 | 1 | 1.00 | U |
| Acenaphthylene | <0.050 | 0.045 | 0.050 | 0.20 | 1 | 1.00 | U |
| Acenaphthene | <0.050 | 0.027 | 0.050 | 0.20 | 1 | 1.00 | U |
| Fluorene | <0.050 | 0.043 | 0.050 | 0.20 | 1 | 1.00 | U |
| Phenanthrene | <0.050 | 0.027 | 0.050 | 0.20 | 1 | 1.00 | U |
| Anthracene | <0.050 | 0.029 | 0.050 | 0.20 | 1 | 1.00 | U |
| Fluoranthene | <0.050 | 0.047 | 0.050 | 0.20 | 1 | 1.00 | U |
| Pyrene | <0.050 | 0.020 | 0.050 | 0.20 | 1 | 1.00 | U |
| Benzo (a) Anthracene | <0.050 | 0.033 | 0.050 | 0.20 | 1 | 1.00 | U |
| Chrysene | <0.050 | 0.025 | 0.050 | 0.20 | 1 | 1.00 | U |
| Benzo (k) Fluoranthene | <0.050 | 0.031 | 0.050 | 0.20 | 1 | 1.00 | U |
| Benzo (b) Fluoranthene | <0.050 | 0.018 | 0.050 | 0.20 | 1 | 1.00 | U |
| Benzo (a) Pyrene | <0.050 | 0.022 | 0.050 | 0.20 | 1 | 1.00 | U |
| Indeno (1,2,3-c,d) Pyrene | <0.050 | 0.021 | 0.050 | 0.20 | 1 | 1.00 | U |
| Dibenz (a,h) Anthracene | <0.050 | 0.048 | 0.050 | 0.20 | 1 | 1.00 | U |
| Benzo (g,h,i) Perylene | <0.10 | 0.082 | 0.10 | 0.20 | . 1 | 1.00 | U |
| <u>Surrogate</u> | Rec. (%) | Control Limi | <u>its Qualifi</u> | <u>ers</u> | | | |
| Nitrobenzene-d5 | 53 | 28-139 | | | | | |
| 2-Fluorobiphenyl | 68 | 33-144 | | | | | |
| p-Terphenyl-d14 | 72 | 23-160 | | | | | |
| | | | | | | | |



Environmental Science International, Inc.

354 Uluniu Street, Suite 304

Preparation: Kailua, HI 96734-2500

Method: Units:

Date Received: 10/28/14 Work Order: 14-10-2131

GC/MS / EPA 8260B

ug/L

EPA 5030C

Project: Red Hill LTM 112066

Page 1 of 12

| Comment(s): | Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|---|-----------------------------------|--------------------------|------------------------|----------------|------------------|-------------------|-----------------------|-------------|
| Parameter Result DL LOD LOQ DE Qualifiers Acetone <10 6.0 10 20 1.00 UICH Bennache <0.50 0.14 0.50 1.0 1.00 UICH Bromodichloromethane <0.50 0.21 0.50 5.0 1.00 UICH Bromomethane <5.0 3.9 5.0 20 1.00 UICH Carbon Tetrachloride <5.0 2.2 5.0 10 1.00 UICH Carbon Tetrachloride <0.50 0.23 0.50 1.0 1.00 UICH Carbon Tetrachloride <0.50 0.23 0.50 1.0 1.00 UICH Carbon Tetrachloride <0.50 0.23 0.50 1.0 1.00 UICH Chlorocethane <0.50 0.17 0.50 5.0 1.00 U Chlorocethane <0.50 0.28 0.50 1.0 1.00 U Li-Diblorocethane <th>ES TRIP</th> <th>14-10-2131-1-A</th> <th></th> <th>Aqueous</th> <th>GC/MS OO</th> <th>10/28/14</th> <th></th> <th>141028L046</th> | ES TRIP | 14-10-2131-1-A | | Aqueous | GC/MS OO | 10/28/14 | | 141028L046 |
| Acetone <10 6.0 10 20 1.00 U.ICH Benzene <0.50 | Comment(s): - Results were evalua | ted to the MDL (DL), con | centrations >= t | to the MDL (DI | L) but < RL (LOC | Q), if found, are | qualified with | a "J" flag. |
| Benzene <0.50 0.14 0.50 1.0 1.00 U Bromodichloromethane <0.50 | <u>Parameter</u> | Result | <u>DL</u> | LOD | LOC | 2 | <u>DF</u> | Qualifiers |
| Bromodichloromethane <0.50 | Acetone | <10 | 6.0 | 10 | 20 | | 1.00 | U,ICH |
| Bromonform <1.0 | Benzene | <0.50 | 0.14 | 0.50 | 1.0 | | 1.00 | U |
| Bromomethane <5.0 | Bromodichloromethane | <0.50 | 0.21 | 0.50 | 5.0 | | 1.00 | U |
| 2-Butanone <5.0 | Bromoform | <1.0 | 0.50 | 1.0 | 10 | | 1.00 | U |
| Carbon Tetrachloride <0.50 | Bromomethane | <5.0 | 3.9 | 5.0 | 20 | | 1.00 | U |
| Chlorobenzene <0.50 0.17 0.50 5.0 1.00 U Chloroethane <5.0 | 2-Butanone | <5.0 | 2.2 | 5.0 | 10 | , | 1.00 | U,ICH |
| Chloroethane <5.0 2.3 5.0 10 1.00 U Chloroform <0.50 | Carbon Tetrachloride | <0.50 | 0.23 | 0.50 | 1.0 | , | 1.00 | U |
| Chloroform < 0.50 0.46 0.50 5.0 1.00 U Chloromethane < 2.0 | Chlorobenzene | <0.50 | 0.17 | 0.50 | 5.0 | | 1.00 | U |
| Chloromethane <2.0 1.8 2.0 10 1.00 U,IJ Dibromochloromethane <0.50 | Chloroethane | <5.0 | 2.3 | 5.0 | 10 | | 1.00 | U |
| Dibromochloromethane 40,50 0,25 0,50 1,0 1,00 U 1,2-Dibromo-3-Chloropropane 42,0 1,2 2,0 10 1,00 U 1,2-Dibromochlane 40,50 0,36 0,50 1,0 1,00 U 1,2-Dibromochlane 40,50 0,46 0,50 1,0 1,00 U 1,3-Dichlorobenzene 40,50 0,40 0,50 1,0 1,00 U 1,4-Dichlorobenzene 40,50 0,43 0,50 1,0 1,00 U 1,1-Dichloroethane 40,50 0,28 0,50 5,0 1,0 1,00 U 1,1-Dichloroethane 40,50 0,28 0,50 5,0 1,0 1,00 U 1,1-Dichloroethane 40,50 0,24 0,50 1,0 1,00 U 1,1-Dichloroethene 40,50 0,43 0,50 1,0 1,00 U 1,1-Dichloroethene 40,50 0,48 0,50 1,0 1,00 U 1,1-Dichloroethene 40,50 0,48 0,50 1,0 1,00 U 1,2-Dichloroethene 40,50 0,48 0,50 1,0 1,00 U 1,2-Dichloroptopane 40,50 0,42 0,50 1,0 1,00 U 1,2-Dichloroptopane 40,50 0,42 0,50 5,0 1,0 1,00 U 1,1-Dichloroptopane 40,50 0,42 0,50 5,0 1,0 1,00 U 1,1-Dichloroptopane 40,50 0,25 0,50 1,0 1,00 U 1,1-Dichloroptopane 40,50 0,25 0,50 1,0 1,00 U 1,1-Dichloroptopane 40,50 0,25 0,50 1,0 1,00 U 4-Methyl-2-Pentanone 40,50 0,14 0,50 1,0 1,00 U 4-Methyl-2-Pentanone 40,50 0,41 0,50 1,0 1,00 U 4-Methyl-2-Pentanone 40,50 0,41 0,50 1,0 1,00 U 1,1,2-Tetrachloroethane 40,50 0,41 0,50 1,0 1,00 U 1,1,1-Trichloroethane 40,50 0,30 0,50 5,0 1,0 1,00 U | Chloroform | <0.50 | 0.46 | 0.50 | 5.0 | , | 1.00 | U |
| 1,2-Dibromo-3-Chloropropane <2.0 | Chloromethane | <2.0 | 1.8 | 2.0 | 10 | | 1.00 | U,IJ |
| 1,2-Dibromoethane <0.50 | Dibromochloromethane | <0.50 | 0.25 | 0.50 | 1.0 | , | 1.00 | U |
| 1,2-Dichlorobenzene <0.50 | 1,2-Dibromo-3-Chloropropane | <2.0 | 1.2 | 2.0 | 10 | | 1.00 | U |
| 1,3-Dichlorobenzene <0.50 | 1,2-Dibromoethane | <0.50 | 0.36 | 0.50 | 1.0 | | 1.00 | U |
| 1,4-Dichlorobenzene <0.50 | 1,2-Dichlorobenzene | <0.50 | 0.46 | 0.50 | 1.0 | | 1.00 | U |
| 1,1-Dichloroethane <0.50 | 1,3-Dichlorobenzene | <0.50 | 0.40 | 0.50 | 1.0 | | 1.00 | U |
| 1,2-Dichloroethane <0.50 | 1,4-Dichlorobenzene | <0.50 | 0.43 | 0.50 | 1.0 | | 1.00 | U |
| 1,1-Dichloroethene <0.50 | 1,1-Dichloroethane | <0.50 | 0.28 | 0.50 | 5.0 | | 1.00 | U |
| c-1,2-Dichloroethene <0.50 | 1,2-Dichloroethane | <0.50 | 0.24 | 0.50 | 1.0 | | 1.00 | U,IH |
| t-1,2-Dichloroethene | 1,1-Dichloroethene | <0.50 | 0.43 | 0.50 | 1.0 | | 1.00 | U |
| 1,2-Dichloropropane <0.50 | c-1,2-Dichloroethene | <0.50 | 0.48 | 0.50 | 1.0 | | 1.00 | U |
| c-1,3-Dichloropropene <0.50 | t-1,2-Dichloroethene | <0.50 | 0.37 | 0.50 | 1.0 | | 1.00 | U |
| t-1,3-Dichloropropene < 0.50 | 1,2-Dichloropropane | <0.50 | 0.42 | 0.50 | 5.0 | | 1.00 | U |
| Ethylbenzene <0.50 0.14 0.50 1.0 1.00 U Methylene Chloride <1.0 | c-1,3-Dichloropropene | <0.50 | 0.25 | 0.50 | 1.0 | | 1.00 | U |
| Methylene Chloride <1.0 0.64 1.0 5.0 1.00 U 4-Methyl-2-Pentanone <5.0 | t-1,3-Dichloropropene | <0.50 | 0.25 | 0.50 | 1.0 | | 1.00 | U |
| 4-Methyl-2-Pentanone <5.0 | Ethylbenzene | <0.50 | 0.14 | 0.50 | 1.0 | | 1.00 | U |
| Styrene <0.50 0.17 0.50 1.0 1.00 U 1,1,1,2-Tetrachloroethane <0.50 | Methylene Chloride | <1.0 | 0.64 | 1.0 | 5.0 | | 1.00 | U |
| 1,1,1,2-Tetrachloroethane <0.50 | 4-Methyl-2-Pentanone | <5.0 | 4.4 | 5.0 | 10 | | 1.00 | U |
| 1,1,1,2-Tetrachloroethane <0.50 | Styrene | <0.50 | 0.17 | 0.50 | 1.0 | | 1.00 | U |
| 1,1,2,2-Tetrachloroethane <0.50 | 1,1,1,2-Tetrachloroethane | | | | | | | U |
| Tetrachloroethene <0.50 0.39 0.50 5.0 1.00 U Toluene <0.50 | 1,1,2,2-Tetrachloroethane | | 0.41 | | | | | |
| Toluene <0.50 0.24 0.50 1.0 1.00 U 1,2,4-Trichlorobenzene <1.0 | Tetrachloroethene | | | | 5.0 | | 1.00 | U |
| 1,2,4-Trichlorobenzene <1.0 | Toluene | | | | | | | |
| 1,1,1-Trichloroethane <0.50 0.30 0.50 5.0 1.00 U | 1,2,4-Trichlorobenzene | | | | | | | |
| | 1,1,1-Trichloroethane | | | | | | | |
| | Hexachloro-1,3-Butadiene | | | | | | | |



Environmental Science International, Inc. 354 Uluniu Street, Suite 304

Kailua, HI 96734-2500

Date Received: Work Order: Preparation:

Method: Units: 10/28/14 14-10-2131

EPA 5030C GC/MS / EPA 8260B

ug/L

Project: Red Hill LTM 112066 Page 2 of 12

| <u>Parameter</u> | Result | <u>DL</u> | LOD | LOQ | <u>DF</u> | Qualifiers |
|-----------------------------|----------|----------------|------------|-----|-----------|------------|
| 1,1,2-Trichloroethane | <0.50 | 0.38 | 0.50 | 1.0 | 1.00 | U |
| Trichloroethene | <0.50 | 0.37 | 0.50 | 1.0 | 1.00 | U |
| 1,2,3-Trichloropropane | <1.0 | 0.64 | 1.0 | 5.0 | 1.00 | U |
| Vinyl Chloride | <0.50 | 0.30 | 0.50 | 1.0 | 1.00 | U |
| p/m-Xylene | <1.0 | 0.30 | 1.0 | 10 | 1.00 | U |
| o-Xylene | <0.50 | 0.23 | 0.50 | 1.0 | 1.00 | U |
| Methyl-t-Butyl Ether (MTBE) | <0.50 | 0.31 | 0.50 | 1.0 | 1.00 | U |
| Gasoline Range Organics | <30 | 26 | 30 | 50 | 1.00 | U |
| Surrogate | Rec. (%) | Control Limits | Qualifiers | | | |
| Dibromofluoromethane | 99 | 80-126 | | | | |
| 1,2-Dichloroethane-d4 | 101 | 80-134 | | | | |
| Toluene-d8 | 100 | 80-120 | | | | |
| Toluene-d8-TPPH | 99 | 88-112 | | | | |
| 1,4-Bromofluorobenzene | 94 | 80-120 | | | | |



Environmental Science International, Inc.

354 Uluniu Street, Suite 304 Kailua, HI 96734-2500 Date Received:

10/28/14 14-10-2131

Work Order: Preparation:

EPA 5030C

Method:

GC/MS / EPA 8260B

Units:

ug/L

Project: Red Hill LTM 112066

Page 3 of 12

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------------------|---------------------------|------------------------|---------------|------------------|-------------------|-----------------------|-------------|
| ES113 | 14-10-2131-2-A | 10/27/14 14:00 | Aqueous | GC/MS OO | 10/28/14 | 10/29/14 06:18 | 141028L046 |
| Comment(s): - Results were evalu | ated to the MDL (DL), con | centrations >= t | o the MDL (DI | _) but < RL (LOC | Q), if found, are | qualified with | a "J" flag. |
| <u>Parameter</u> | Result | <u>DL</u> | <u>LOD</u> | LOC | <u>D</u> | <u>E</u> | Qualifiers |
| Acetone | <10 | 6.0 | 10 | 20 | 1 | .00 | U,ICH |
| Benzene | <0.50 | 0.14 | 0.50 | 1.0 | 1 | .00 | U |
| Bromodichloromethane | <0.50 | 0.21 | 0.50 | 5.0 | 1 | .00 | U |
| Bromoform | <1.0 | 0.50 | 1.0 | 10 | 1 | .00 | U |
| Bromomethane | <5.0 | 3.9 | 5.0 | 20 | 1 | .00 | U |
| 2-Butanone | <5.0 | 2.2 | 5.0 | 10 | 1 | .00 | U,ICH |
| Carbon Tetrachloride | <0.50 | 0.23 | 0.50 | 1.0 | 1 | .00 | U |
| Chlorobenzene | <0.50 | 0.17 | 0.50 | 5.0 | 1 | .00 | U |
| Chloroethane | <5.0 | 2.3 | 5.0 | 10 | 1 | .00 | U |
| Chloroform | <0.50 | 0.46 | 0.50 | 5.0 | 1 | .00 | U |
| Chloromethane | <2.0 | 1.8 | 2.0 | 10 | 1 | .00 | U,IJ |
| Dibromochloromethane | <0.50 | 0.25 | 0.50 | 1.0 | 1 | .00 | U |
| 1,2-Dibromo-3-Chloropropane | <2.0 | 1.2 | 2.0 | 10 | 1 | .00 | U |
| 1,2-Dibromoethane | <0.50 | 0.36 | 0.50 | 1.0 | 1 | .00 | U |
| 1,2-Dichlorobenzene | <0.50 | 0.46 | 0.50 | 1.0 | 1 | .00 | U |
| 1,3-Dichlorobenzene | <0.50 | 0.40 | 0.50 | 1.0 | 1 | .00 | U |
| 1,4-Dichlorobenzene | <0.50 | 0.43 | 0.50 | 1.0 | 1 | .00 | U |
| 1,1-Dichloroethane | <0.50 | 0.28 | 0.50 | 5.0 | 1 | .00 | U |
| 1,2-Dichloroethane | <0.50 | 0.24 | 0.50 | 1.0 | 1 | .00 | U,IH |
| 1,1-Dichloroethene | <0.50 | 0.43 | 0.50 | 1.0 | 1 | .00 | U |
| c-1,2-Dichloroethene | <0.50 | 0.48 | 0.50 | 1.0 | 1 | .00 | U |
| t-1,2-Dichloroethene | <0.50 | 0.37 | 0.50 | 1.0 | 1 | .00 | U |
| 1,2-Dichloropropane | <0.50 | 0.42 | 0.50 | 5.0 | 1 | .00 | U |
| c-1,3-Dichloropropene | <0.50 | 0.25 | 0.50 | 1.0 | 1 | .00 | U |
| t-1,3-Dichloropropene | <0.50 | 0.25 | 0.50 | 1.0 | 1 | .00 | U |
| Ethylbenzene | <0.50 | 0.14 | 0.50 | 1.0 | 1 | .00 | U |
| Methylene Chloride | <1.0 | 0.64 | 1.0 | 5.0 | 1 | .00 | U |
| 4-Methyl-2-Pentanone | <5.0 | 4.4 | 5.0 | 10 | 1 | .00 | U |
| Styrene | <0.50 | 0.17 | 0.50 | 1.0 | 1 | .00 | U |
| 1,1,1,2-Tetrachloroethane | <0.50 | 0.40 | 0.50 | 1.0 | | .00 | U |
| 1,1,2,2-Tetrachloroethane | <0.50 | 0.41 | 0.50 | 1.0 | 1 | .00 | U |
| Tetrachloroethene | <0.50 | 0.39 | 0.50 | 5.0 | 1 | .00 | U |
| Toluene | <0.50 | 0.24 | 0.50 | 1.0 | 1 | .00 | U |
| 1,2,4-Trichlorobenzene | <1.0 | 0.50 | 1.0 | 5.0 | 1 | .00 | U |
| 1,1,1-Trichloroethane | <0.50 | 0.30 | 0.50 | 5.0 | 1 | .00 | U |
| Hexachloro-1,3-Butadiene | <0.50 | 0.32 | 0.50 | 1.0 | 1 | .00 | U |



Environmental Science International, Inc. 354 Uluniu Street, Suite 304

Kailua, HI 96734-2500

Date Received:

Work Order:

Preparation:

Method: Units: 10/28/14 14-10-2131

EPA 5030C

GC/MS / EPA 8260B

ug/L

Project: Red Hill LTM 112066

Page 4 of 12

| <u>Parameter</u> | <u>Result</u> | <u>DL</u> | LOD | LOQ | <u>DF</u> | <u>Qualifiers</u> |
|-----------------------------|---------------|----------------|------------|-----|-----------|-------------------|
| 1,1,2-Trichloroethane | <0.50 | 0.38 | 0.50 | 1.0 | 1.00 | U |
| Trichloroethene | <0.50 | 0.37 | 0.50 | 1.0 | 1.00 | U |
| 1,2,3-Trichloropropane | <1.0 | 0.64 | 1.0 | 5.0 | 1.00 | U |
| Vinyl Chloride | <0.50 | 0.30 | 0.50 | 1.0 | 1.00 | U |
| p/m-Xylene | <1.0 | 0.30 | 1.0 | 10 | 1.00 | U |
| o-Xylene | <0.50 | 0.23 | 0.50 | 1.0 | 1.00 | U |
| Methyl-t-Butyl Ether (MTBE) | <0.50 | 0.31 | 0.50 | 1.0 | 1.00 | U |
| Gasoline Range Organics | <30 | 26 | 30 | 50 | 1.00 | U |
| Surrogate | Rec. (%) | Control Limits | Qualifiers | | | |
| Dibromofluoromethane | 95 | 80-126 | | | | |
| 1,2-Dichloroethane-d4 | 97 | 80-134 | | | | |
| Toluene-d8 | 99 | 80-120 | | | | |
| Toluene-d8-TPPH | 98 | 88-112 | | | | |
| 1,4-Bromofluorobenzene | 95 | 80-120 | | | | |

10/28/14



Analytical Report

Date Received: Environmental Science International, Inc. 354 Uluniu Street, Suite 304 Work Order: 14-10-2131

Preparation: EPA 5030C Kailua, HI 96734-2500 Method: GC/MS / EPA 8260B

> Units: ug/L

Project: Red Hill LTM 112066 Page 5 of 12

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|---|----------------------|------------------------|---------------|------------------|------------------|-----------------------|-------------------|
| ES114 | 14-10-2131-3-A | 10/27/14 10:45 | Aqueous | GC/MS OO | 10/28/14 | 10/29/14 05:51 | 141028L046 |
| Comment(s): - Results were evaluated to | o the MDL (DL), con- | centrations >= t | o the MDL (DI | L) but < RL (LOQ |), if found, a | re qualified with a | a "J" flag. |
| <u>Parameter</u> | <u>Result</u> | <u>DL</u> | <u>LOD</u> | LOC | <u>!</u> | <u>DF</u> | <u>Qualifiers</u> |
| Acetone | <10 | 6.0 | 10 | 20 | | 1.00 | U,ICH |
| Benzene | <0.50 | 0.14 | 0.50 | 1.0 | | 1.00 | U |
| Bromodichloromethane | <0.50 | 0.21 | 0.50 | 5.0 | | 1.00 | U |
| Bromoform | <1.0 | 0.50 | 1.0 | 10 | | 1.00 | U |
| Bromomethane | <5.0 | 3.9 | 5.0 | 20 | | 1.00 | U |
| 2-Butanone | <5.0 | 2.2 | 5.0 | 10 | | 1.00 | U,ICH |
| Carbon Tetrachloride | <0.50 | 0.23 | 0.50 | 1.0 | | 1.00 | U |
| Chlorobenzene | <0.50 | 0.17 | 0.50 | 5.0 | | 1.00 | U |
| Chloroethane | <5.0 | 2.3 | 5.0 | 10 | | 1.00 | U |
| Chloroform | <0.50 | 0.46 | 0.50 | 5.0 | | 1.00 | U |
| Chloromethane | <2.0 | 1.8 | 2.0 | 10 | | 1.00 | U,IJ |
| Dibromochloromethane | <0.50 | 0.25 | 0.50 | 1.0 | | 1.00 | U |
| 1,2-Dibromo-3-Chloropropane | <2.0 | 1.2 | 2.0 | 10 | | 1.00 | U |
| 1,2-Dibromoethane | <0.50 | 0.36 | 0.50 | 1.0 | | 1.00 | U |
| 1,2-Dichlorobenzene | <0.50 | 0.46 | 0.50 | 1.0 | | 1.00 | U |
| 1,3-Dichlorobenzene | <0.50 | 0.40 | 0.50 | 1.0 | | 1.00 | U |
| 1,4-Dichlorobenzene | <0.50 | 0.43 | 0.50 | 1.0 | | 1.00 | U |
| 1,1-Dichloroethane | <0.50 | 0.28 | 0.50 | 5.0 | | 1.00 | U |
| 1,2-Dichloroethane | <0.50 | 0.24 | 0.50 | 1.0 | | 1.00 | U,IH |
| 1,1-Dichloroethene | <0.50 | 0.43 | 0.50 | 1.0 | | 1.00 | U |
| c-1,2-Dichloroethene | <0.50 | 0.48 | 0.50 | 1.0 | | 1.00 | U |
| t-1,2-Dichloroethene | <0.50 | 0.37 | 0.50 | 1.0 | | 1.00 | U |
| 1,2-Dichloropropane | <0.50 | 0.42 | 0.50 | 5.0 | | 1.00 | U |
| c-1,3-Dichloropropene | <0.50 | 0.25 | 0.50 | 1.0 | | 1.00 | U |
| t-1,3-Dichloropropene | <0.50 | 0.25 | 0.50 | 1.0 | | 1.00 | U |
| Ethylbenzene | 0.15 | 0.14 | 0.50 | 1.0 | | 1.00 | J |
| Methylene Chloride | <1.0 | 0.64 | 1.0 | 5.0 | | 1.00 | U |
| 4-Methyl-2-Pentanone | <5.0 | 4.4 | 5.0 | 10 | | 1.00 | U |
| Styrene | <0.50 | 0.17 | 0.50 | 1.0 | | 1.00 | U |
| 1,1,1,2-Tetrachloroethane | <0.50 | 0.40 | 0.50 | 1.0 | | 1.00 | U |
| 1,1,2,2-Tetrachloroethane | <0.50 | 0.41 | 0.50 | 1.0 | | 1.00 | U |
| Tetrachloroethene | <0.50 | 0.39 | 0.50 | 5.0 | | 1.00 | U |
| Toluene | <0.50 | 0.24 | 0.50 | 1.0 | | 1.00 | U |
| 1,2,4-Trichlorobenzene | <1.0 | 0.50 | 1.0 | 5.0 | | 1.00 | U |
| 1,1,1-Trichloroethane | <0.50 | 0.30 | 0.50 | 5.0 | | 1.00 | U |
| Hexachloro-1,3-Butadiene | <0.50 | 0.32 | 0.50 | 1.0 | | 1.00 | U |



Environmental Science International, Inc. 354 Uluniu Street, Suite 304

Kailua, HI 96734-2500

Date Received: Work Order:

Preparation:

Method: Units: 10/28/14 14-10-2131

EPA 5030C

GC/MS / EPA 8260B

ug/

Project: Red Hill LTM 112066

Page 6 of 12

| Parameter | <u>Result</u> | <u>DL</u> | LOD | LOQ | <u>DF</u> | Qualifiers |
|-----------------------------|---------------|----------------|------------|-----|-----------|------------|
| 1,1,2-Trichloroethane | <0.50 | 0.38 | 0.50 | 1.0 | 1.00 | U |
| Trichloroethene | <0.50 | 0.37 | 0.50 | 1.0 | 1.00 | U |
| 1,2,3-Trichloropropane | <1.0 | 0.64 | 1.0 | 5.0 | 1.00 | U |
| Vinyl Chloride | <0.50 | 0.30 | 0.50 | 1.0 | 1.00 | U |
| p/m-Xylene | <1.0 | 0.30 | 1.0 | 10 | 1.00 | U |
| o-Xylene | 0.32 | 0.23 | 0.50 | 1.0 | 1.00 | J |
| Methyl-t-Butyl Ether (MTBE) | <0.50 | 0.31 | 0.50 | 1.0 | 1.00 | U |
| Gasoline Range Organics | 57 | 26 | 30 | 50 | 1.00 | |
| Surrogate | Rec. (%) | Control Limits | Qualifiers | | | |
| Dibromofluoromethane | 98 | 80-126 | | | | |
| 1,2-Dichloroethane-d4 | 100 | 80-134 | | | | |
| Toluene-d8 | 98 | 80-120 | | | | |
| Toluene-d8-TPPH | 97 | 88-112 | | | | |
| 1 4-Bromofluorobenzene | 98 | 80-120 | | | | |



Environmental Science International, Inc.

Date Received: 10/28/14

354 Uluniu Street, Suite 304

Work Order: 14-10-2131

 Kailua, HI 96734-2500
 Preparation:
 EPA 5030C

 Method:
 GC/MS / EPA 8260B

Units: ug/L

Project: Red Hill LTM 112066 Page 7 of 12

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Date/Time Prepared Analyzed | QC Batch ID |
|-------------------------------------|-------------------------|------------------------|---------------|------------------|-------------------------------------|-------------|
| ES115 | 14-10-2131-4-A | 10/27/14 11:15 | Aqueous | GC/MS OO | 10/28/14 10/29/14 06:45 | 141028L046 |
| Comment(s): - Results were evaluate | ed to the MDL (DL), con | centrations >= t | o the MDL (DI | _) but < RL (LOC | t), if found, are qualified with | a "J" flag. |
| Parameter | Result | <u>DL</u> | LOD | LOC | <u>DF</u> | Qualifiers |
| Acetone | <10 | 6.0 | 10 | 20 | 1.00 | U,ICH |
| Benzene | <0.50 | 0.14 | 0.50 | 1.0 | 1.00 | U |
| Bromodichloromethane | <0.50 | 0.21 | 0.50 | 5.0 | 1.00 | U |
| Bromoform | <1.0 | 0.50 | 1.0 | 10 | 1.00 | U |
| Bromomethane | <5.0 | 3.9 | 5.0 | 20 | 1.00 | U |
| 2-Butanone | <5.0 | 2.2 | 5.0 | 10 | 1.00 | U,ICH |
| Carbon Tetrachloride | <0.50 | 0.23 | 0.50 | 1.0 | 1.00 | U |
| Chlorobenzene | <0.50 | 0.17 | 0.50 | 5.0 | 1.00 | U |
| Chloroethane | <5.0 | 2.3 | 5.0 | 10 | 1.00 | U |
| Chloroform | <0.50 | 0.46 | 0.50 | 5.0 | 1.00 | U |
| Chloromethane | <2.0 | 1.8 | 2.0 | 10 | 1.00 | U,IJ |
| Dibromochloromethane | <0.50 | 0.25 | 0.50 | 1.0 | 1.00 | U |
| 1,2-Dibromo-3-Chloropropane | <2.0 | 1.2 | 2.0 | 10 | 1.00 | U |
| 1,2-Dibromoethane | <0.50 | 0.36 | 0.50 | 1.0 | 1.00 | U |
| 1,2-Dichlorobenzene | <0.50 | 0.46 | 0.50 | 1.0 | 1.00 | U |
| 1,3-Dichlorobenzene | <0.50 | 0.40 | 0.50 | 1.0 | 1.00 | U |
| 1,4-Dichlorobenzene | <0.50 | 0.43 | 0.50 | 1.0 | 1.00 | U |
| 1,1-Dichloroethane | <0.50 | 0.28 | 0.50 | 5.0 | 1.00 | U |
| 1,2-Dichloroethane | <0.50 | 0.24 | 0.50 | 1.0 | 1.00 | U,IH |
| 1,1-Dichloroethene | <0.50 | 0.43 | 0.50 | 1.0 | 1.00 | U |
| c-1,2-Dichloroethene | <0.50 | 0.48 | 0.50 | 1.0 | 1.00 | U |
| t-1,2-Dichloroethene | <0.50 | 0.37 | 0.50 | 1.0 | 1.00 | U |
| 1,2-Dichloropropane | <0.50 | 0.42 | 0.50 | 5.0 | 1.00 | U |
| c-1,3-Dichloropropene | <0.50 | 0.25 | 0.50 | 1.0 | 1.00 | U |
| t-1,3-Dichloropropene | <0.50 | 0.25 | 0.50 | 1.0 | 1.00 | U |
| Ethylbenzene | 0.14 | 0.14 | 0.50 | 1.0 | 1.00 | J |
| Methylene Chloride | <1.0 | 0.64 | 1.0 | 5.0 | 1.00 | U |
| 4-Methyl-2-Pentanone | <5.0 | 4.4 | 5.0 | 10 | 1.00 | U |
| Styrene | <0.50 | 0.17 | 0.50 | 1.0 | 1.00 | U |
| 1,1,1,2-Tetrachloroethane | <0.50 | 0.40 | 0.50 | 1.0 | 1.00 | U |
| 1,1,2,2-Tetrachloroethane | <0.50 | 0.41 | 0.50 | 1.0 | 1.00 | U |
| Tetrachloroethene | <0.50 | 0.39 | 0.50 | 5.0 | 1.00 | U |
| Toluene | <0.50 | 0.24 | 0.50 | 1.0 | 1.00 | U |
| 1,2,4-Trichlorobenzene | <1.0 | 0.50 | 1.0 | 5.0 | 1.00 | U |
| 1,1,1-Trichloroethane | <0.50 | 0.30 | 0.50 | 5.0 | 1.00 | U |
| Hexachloro-1,3-Butadiene | <0.50 | 0.32 | 0.50 | 1.0 | 1.00 | U |



Environmental Science International, Inc. 354 Uluniu Street, Suite 304

Kailua, HI 96734-2500

Date Received: Work Order:

Preparation: Method: Units: 10/28/14 14-10-2131

EPA 5030C

GC/MS / EPA 8260B ug/L

Page 8 of 12

Qualifiers
U
U
U
U
U
J

| Project: Red Hill LTM 112066 | | | | | |
|------------------------------|---------------|----------------|------------|-----|-----------|
| <u>Parameter</u> | <u>Result</u> | <u>DL</u> | LOD | LOQ | <u>DF</u> |
| 1,1,2-Trichloroethane | <0.50 | 0.38 | 0.50 | 1.0 | 1.00 |
| Trichloroethene | <0.50 | 0.37 | 0.50 | 1.0 | 1.00 |
| 1,2,3-Trichloropropane | <1.0 | 0.64 | 1.0 | 5.0 | 1.00 |
| Vinyl Chloride | <0.50 | 0.30 | 0.50 | 1.0 | 1.00 |
| p/m-Xylene | <1.0 | 0.30 | 1.0 | 10 | 1.00 |
| o-Xylene | 0.29 | 0.23 | 0.50 | 1.0 | 1.00 |
| Methyl-t-Butyl Ether (MTBE) | <0.50 | 0.31 | 0.50 | 1.0 | 1.00 |
| Gasoline Range Organics | 53 | 26 | 30 | 50 | 1.00 |
| Surrogate | Rec. (%) | Control Limits | Qualifiers | | |
| Dibromofluoromethane | 96 | 80-126 | | | |
| 1,2-Dichloroethane-d4 | 97 | 80-134 | | | |
| Toluene-d8 | 99 | 80-120 | | | |
| Toluene-d8-TPPH | 98 | 88-112 | | | |
| 1,4-Bromofluorobenzene | 96 | 80-120 | | | |

10/28/14

14-10-2131



Analytical Report

Environmental Science International, Inc.

Date Received:

Work Order:

 Kailua, HI 96734-2500
 Preparation:
 EPA 5030C

 Method:
 GC/MS / EPA 8260B

Units: ug/L

Project: Red Hill LTM 112066 Page 9 of 12

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Date/Time Prepared Analyzed | QC Batch ID |
|--------------------------------------|-----------------------|------------------------|---------------|------------------|-------------------------------------|-------------------|
| ES116 | 14-10-2131-5-A | 10/27/14 09:00 | Aqueous | GC/MS OO | 10/28/14 10/29/14 07:13 | 141028L046 |
| Comment(s): - Results were evaluated | to the MDL (DL), con- | centrations >= t | o the MDL (DL | _) but < RL (LOC |), if found, are qualified with | n a "J" flag. |
| Parameter | <u>Result</u> | <u>DL</u> | LOD | LOG | <u>DF</u> | <u>Qualifiers</u> |
| Acetone | <10 | 6.0 | 10 | 20 | 1.00 | U,ICH |
| Benzene | <0.50 | 0.14 | 0.50 | 1.0 | 1.00 | U |
| Bromodichloromethane | <0.50 | 0.21 | 0.50 | 5.0 | 1.00 | U |
| Bromoform | <1.0 | 0.50 | 1.0 | 10 | 1.00 | U |
| Bromomethane | <5.0 | 3.9 | 5.0 | 20 | 1.00 | U |
| 2-Butanone | <5.0 | 2.2 | 5.0 | 10 | 1.00 | U,ICH |
| Carbon Tetrachloride | <0.50 | 0.23 | 0.50 | 1.0 | 1.00 | U |
| Chlorobenzene | <0.50 | 0.17 | 0.50 | 5.0 | 1.00 | U |
| Chloroethane | <5.0 | 2.3 | 5.0 | 10 | 1.00 | U |
| Chloroform | <0.50 | 0.46 | 0.50 | 5.0 | 1.00 | U |
| Chloromethane | <2.0 | 1.8 | 2.0 | 10 | 1.00 | U,IJ |
| Dibromochloromethane | <0.50 | 0.25 | 0.50 | 1.0 | 1.00 | U |
| 1,2-Dibromo-3-Chloropropane | <2.0 | 1.2 | 2.0 | 10 | 1.00 | U |
| 1,2-Dibromoethane | <0.50 | 0.36 | 0.50 | 1.0 | 1.00 | U |
| 1,2-Dichlorobenzene | <0.50 | 0.46 | 0.50 | 1.0 | 1.00 | U |
| 1,3-Dichlorobenzene | <0.50 | 0.40 | 0.50 | 1.0 | 1.00 | U |
| 1,4-Dichlorobenzene | <0.50 | 0.43 | 0.50 | 1.0 | 1.00 | U |
| 1,1-Dichloroethane | <0.50 | 0.28 | 0.50 | 5.0 | 1.00 | U |
| 1,2-Dichloroethane | <0.50 | 0.24 | 0.50 | 1.0 | 1.00 | U,IH |
| 1,1-Dichloroethene | <0.50 | 0.43 | 0.50 | 1.0 | 1.00 | U |
| c-1,2-Dichloroethene | <0.50 | 0.48 | 0.50 | 1.0 | 1.00 | U |
| t-1,2-Dichloroethene | <0.50 | 0.37 | 0.50 | 1.0 | 1.00 | U |
| 1,2-Dichloropropane | <0.50 | 0.42 | 0.50 | 5.0 | 1.00 | U |
| c-1,3-Dichloropropene | <0.50 | 0.25 | 0.50 | 1.0 | 1.00 | U |
| t-1,3-Dichloropropene | <0.50 | 0.25 | 0.50 | 1.0 | 1.00 | U |
| Ethylbenzene | <0.50 | 0.14 | 0.50 | 1.0 | 1.00 | U |
| Methylene Chloride | <1.0 | 0.64 | 1.0 | 5.0 | 1.00 | U |
| 4-Methyl-2-Pentanone | <5.0 | 4.4 | 5.0 | 10 | 1.00 | U |
| Styrene | <0.50 | 0.17 | 0.50 | 1.0 | 1.00 | U |
| 1,1,1,2-Tetrachloroethane | <0.50 | 0.40 | 0.50 | 1.0 | 1.00 | U |
| 1,1,2,2-Tetrachloroethane | <0.50 | 0.41 | 0.50 | 1.0 | 1.00 | U |
| Tetrachloroethene | <0.50 | 0.39 | 0.50 | 5.0 | 1.00 | U |
| Toluene | <0.50 | 0.24 | 0.50 | 1.0 | 1.00 | U |
| 1,2,4-Trichlorobenzene | <1.0 | 0.50 | 1.0 | 5.0 | 1.00 | U |
| 1,1,1-Trichloroethane | <0.50 | 0.30 | 0.50 | 5.0 | 1.00 | U |
| Hexachloro-1,3-Butadiene | <0.50 | 0.32 | 0.50 | 1.0 | 1.00 | U |



Environmental Science International, Inc. 354 Uluniu Street, Suite 304 Kailua, HI 96734-2500 Date Received:
Work Order:
Preparation:
Method:

10/28/14 14-10-2131 EPA 5030C GC/MS / EPA 8260B

ug/L

Units:

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| Proje | oct. | Red | Hill | I TN/I | 11 | 2066 |
|-------|------|-----|--------|---------|-----|------|
| | 7UL. | Neu | 1 1111 | L I IVI | 3 1 | 2000 |

| <u>Parameter</u> | Result | <u>DL</u> | LOD | LOQ | <u>DF</u> | Qualifiers |
|-----------------------------|----------|----------------|------------|-----|-----------|------------|
| 1,1,2-Trichloroethane | <0.50 | 0.38 | 0.50 | 1.0 | 1.00 | U |
| Trichloroethene | <0.50 | 0.37 | 0.50 | 1.0 | 1.00 | U |
| 1,2,3-Trichloropropane | <1.0 | 0.64 | 1.0 | 5.0 | 1.00 | U |
| Vinyl Chloride | <0.50 | 0.30 | 0.50 | 1.0 | 1.00 | U |
| p/m-Xylene | <1.0 | 0.30 | 1.0 | 10 | 1.00 | U |
| o-Xylene | <0.50 | 0.23 | 0.50 | 1.0 | 1.00 | U |
| Methyl-t-Butyl Ether (MTBE) | <0.50 | 0.31 | 0.50 | 1.0 | 1.00 | U |
| Gasoline Range Organics | <30 | 26 | 30 | 50 | 1.00 | U |
| Surrogate | Rec. (%) | Control Limits | Qualifiers | | | |
| Dibromofluoromethane | 95 | 80-126 | | | | |
| 1,2-Dichloroethane-d4 | 98 | 80-134 | | | | |
| Toluene-d8 | 99 | 80-120 | | | | |
| Toluene-d8-TPPH | 98 | 88-112 | | | | |
| 1,4-Bromofluorobenzene | 95 | 80-120 | | | | |



Environmental Science International, Inc.

354 Uluniu Street, Suite 304

Kailua, HI 96734-2500

Date Received:

10/28/14 Work Order: 14-10-2131

Preparation:

EPA 5030C

Method: Units:

GC/MS / EPA 8260B

ug/L Page 11 of 12

Project: Red Hill LTM 112066

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------------------|---------------------------|------------------------|---------------|------------------|-------------------|-----------------------|-------------------|
| Method Blank | 099-13-057-71 | N/A | Aqueous | GC/MS OO | 10/28/14 | 10/29/14 04:56 | 141028L046 |
| Comment(s): - Results were evalu | ated to the MDL (DL), cor | ncentrations >= t | o the MDL (DI | _) but < RL (LOC | Q), if found, are | qualified with | a "J" flag. |
| <u>Parameter</u> | <u>Result</u> | <u>DL</u> | LOD | LOC | <u> </u> | <u>DF</u> | <u>Qualifiers</u> |
| Acetone | <10 | 6.0 | 10 | 20 | | 1.00 | U |
| Benzene | <0.50 | 0.14 | 0.50 | 1.0 | | 1.00 | U |
| Bromodichloromethane | <0.50 | 0.21 | 0.50 | 5.0 | • | 1.00 | U |
| Bromoform | <1.0 | 0.50 | 1.0 | 10 | | 1.00 | U |
| Bromomethane | <5.0 | 3.9 | 5.0 | 20 | | 1.00 | U |
| 2-Butanone | <5.0 | 2.2 | 5.0 | 10 | | 1.00 | U |
| Carbon Tetrachloride | <0.50 | 0.23 | 0.50 | 1.0 | | 1.00 | U |
| Chlorobenzene | <0.50 | 0.17 | 0.50 | 5.0 | | 1.00 | U |
| Chloroethane | <5.0 | 2.3 | 5.0 | 10 | | 1.00 | U |
| Chloroform | <0.50 | 0.46 | 0.50 | 5.0 | | 1.00 | U |
| Chloromethane | <2.0 | 1.8 | 2.0 | 10 | | 1.00 | U |
| Dibromochloromethane | <0.50 | 0.25 | 0.50 | 1.0 | | 1.00 | U |
| 1,2-Dibromo-3-Chloropropane | <2.0 | 1.2 | 2.0 | 10 | | 1.00 | U |
| 1,2-Dibromoethane | <0.50 | 0.36 | 0.50 | 1.0 | | 1.00 | U |
| 1,2-Dichlorobenzene | <0.50 | 0.46 | 0.50 | 1.0 | | 1.00 | U |
| 1,3-Dichlorobenzene | <0.50 | 0.40 | 0.50 | 1.0 | | 1.00 | U |
| 1,4-Dichlorobenzene | <0.50 | 0.43 | 0.50 | 1.0 | • | 1.00 | U |
| 1,1-Dichloroethane | <0.50 | 0.28 | 0.50 | 5.0 | | 1.00 | U |
| 1,2-Dichloroethane | <0.50 | 0.24 | 0.50 | 1.0 | | 1.00 | U |
| 1,1-Dichloroethene | <0.50 | 0.43 | 0.50 | 1.0 | | 1.00 | U |
| c-1,2-Dichloroethene | <0.50 | 0.48 | 0.50 | 1.0 | | 1.00 | U |
| t-1,2-Dichloroethene | <0.50 | 0.37 | 0.50 | 1.0 | | 1.00 | U |
| 1,2-Dichloropropane | <0.50 | 0.42 | 0.50 | 5.0 | • | 1.00 | U |
| c-1,3-Dichloropropene | <0.50 | 0.25 | 0.50 | 1.0 | | 1.00 | U |
| -1,3-Dichloropropene | <0.50 | 0.25 | 0.50 | 1.0 | | 1.00 | U |
| Ethylbenzene | <0.50 | 0.14 | 0.50 | 1.0 | | 1.00 | U |
| Methylene Chloride | <1.0 | 0.64 | 1.0 | 5.0 | | 1.00 | U |
| 4-Methyl-2-Pentanone | <5.0 | 4.4 | 5.0 | 10 | | 1.00 | U |
| Styrene | <0.50 | 0.17 | 0.50 | 1.0 | | 1.00 | U |
| 1,1,1,2-Tetrachloroethane | <0.50 | 0.40 | 0.50 | 1.0 | | 1.00 | U |
| 1,1,2,2-Tetrachloroethane | <0.50 | 0.41 | 0.50 | 1.0 | | 1.00 | U |
| Tetrachloroethene | <0.50 | 0.39 | 0.50 | 5.0 | | 1.00 | U |
| Toluene | <0.50 | 0.24 | 0.50 | 1.0 | | 1.00 | U |
| 1,2,4-Trichlorobenzene | <1.0 | 0.50 | 1.0 | 5.0 | | 1.00 | U |
| 1,1,1-Trichloroethane | <0.50 | 0.30 | 0.50 | 5.0 | | 1.00 | U |
| Hexachloro-1,3-Butadiene | <0.50 | 0.32 | 0.50 | 1.0 | | 1.00 | U |



Environmental Science International, Inc. 354 Uluniu Street, Suite 304

Kailua, HI 96734-2500

Date Received:

Work Order:

Preparation: Method:

GC/MS / EPA 8260B

Units:

ug/L

10/28/14

14-10-2131

EPA 5030C

Project: Red Hill LTM 112066

Page 12 of 12

| <u>Parameter</u> | Result | <u>DL</u> | LOD | LOQ | DF | <u>Qualifiers</u> |
|-----------------------------|----------|----------------|------------|-----|------|-------------------|
| 1,1,2-Trichloroethane | <0.50 | 0.38 | 0.50 | 1.0 | 1.00 | U |
| Trichloroethene | <0.50 | 0.37 | 0.50 | 1.0 | 1.00 | U |
| 1,2,3-Trichloropropane | <1.0 | 0.64 | 1.0 | 5.0 | 1.00 | U |
| Vinyl Chloride | <0.50 | 0.30 | 0.50 | 1.0 | 1.00 | U |
| p/m-Xylene | <1.0 | 0.30 | 1.0 | 10 | 1.00 | U |
| o-Xylene | <0.50 | 0.23 | 0.50 | 1.0 | 1.00 | U |
| Methyl-t-Butyl Ether (MTBE) | <0.50 | 0.31 | 0.50 | 1.0 | 1.00 | U |
| Gasoline Range Organics | <30 | 26 | 30 | 50 | 1.00 | U |
| Surrogate | Rec. (%) | Control Limits | Qualifiers | | | |
| Dibromofluoromethane | 98 | 80-126 | | | | |
| 1,2-Dichloroethane-d4 | 101 | 80-134 | | | | |
| Toluene-d8 | 100 | 80-120 | | | | |
| Toluene-d8-TPPH | 100 | 88-112 | | | | |
| 1,4-Bromofluorobenzene | 95 | 80-120 | | | | |



Quality Control - Spike/Spike Duplicate

Environmental Science International, Inc.

Date Received:

Work Order:

14-10-2131

Kailua, HI 96734-2500

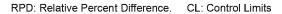
Preparation:

Method:

EPA 8015B (M)

| Project: Red Hill LTM 112066 | Page 1 of 5 |
|------------------------------|-------------|
|------------------------------|-------------|

| Quality Control Sample ID | Туре | Matrix | Instrument | Date Prepared | Date Analyzed | MS/MSD Bat | tch Number |
|---------------------------|--|-------------|--------------------------------------|---------------|----------------|------------|------------|
| ES114 | Sample | Aqueous | GC 45 | 10/30/14 | 10/31/14 09:59 | 141030S06 | |
| ES114 | Matrix Spike | Aqueous | GC 45 | 10/30/14 | 10/31/14 09:05 | 141030S06 | |
| ES114 | Matrix Spike Duplica | ite Aqueous | GC 45 | 10/30/14 | 10/31/14 09:23 | 141030S06 | |
| <u>Parameter</u> | <u>Sample</u> Spike <u>Conc.</u> Adde | | <u>MS</u> <u>MS</u> <u>%Rec.</u> Cor | | %Rec. CL RPD | RPD CL | Qualifiers |
| TPH as Diesel | 2046 2000 | 7915 | 293 799 | 4 297 | 55-133 1 | 0-30 | 3 |





Quality Control - Spike/Spike Duplicate

Environmental Science International, Inc.

Date Received:

Work Order:

14-10-2131

Kailua, HI 96734-2500

Preparation:

Method:

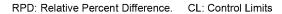
EPA 3005A Filt.

Method:

Project: Red Hill LTM 112066 Page 2 of 5

| Quality Control Sample ID | Туре | | Matrix | Ins | strument | Date Prepared | Date Ana | lyzed | MS/MSD Ba | tch Number |
|---------------------------|-------------------------------|-----------------------|-------------|--------------------|--------------|---------------|----------|-------|------------|------------|
| ES114 | Sample | | Aqueou | s IC | P/MS 04 | 10/29/14 | 10/30/14 | 23:32 | 141029S03E | 3 |
| ES114 | Matrix Spike | | Aqueou | s IC | P/MS 04 | 10/29/14 | 10/30/14 | 23:05 | 141029S03E | 3 |
| ES114 | Matrix Spike E | Ouplicate | Aqueou | s IC | P/MS 04 | 10/29/14 | 10/30/14 | 23:19 | 141029S03E | 3 |
| Parameter | <u>Sample</u> <u>Conc.</u> | <u>Spike</u> Added | MS Conc. | <u>MS</u> %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Lead | ND | 100.0 | 112.2 | 112 | 112.7 | 113 | 80-120 | 0 | 0-20 | |

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EPA 8270C SIM PAHs



Quality Control - Spike/Spike Duplicate

Environmental Science International, Inc.

Date Received:

Work Order:

10/28/14

Work Order:

Kailua, HI 96734-2500

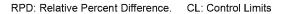
Preparation:

EPA 3510C

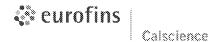
Method:

Project: Red Hill LTM 112066 Page 3 of 5

| Quality Control Sample ID | Туре | | Matrix | Ir | nstrument | Date Prepared | Date Ana | lyzed | MS/MSD Ba | tch Number |
|---------------------------|-------------------------------|------------------------------|-------------|-------------|--------------|---------------|----------|-------|-----------|------------|
| ES114 | Sample | | Aqueou | ıs C | GC/MS AAA | 10/30/14 | 11/03/14 | 18:28 | 141030S08 | |
| ES114 | Matrix Spike | | Aqueou | ıs C | GC/MS AAA | 10/30/14 | 10/31/14 | 18:04 | 141030508 | |
| ES114 | Matrix Spike | Duplicate | Aqueou | ıs C | GC/MS AAA | 10/30/14 | 10/31/14 | 18:27 | 141030508 | |
| <u>Parameter</u> | <u>Sample</u> <u>Conc.</u> | <u>Spike</u> <u>Added</u> | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Naphthalene | 143.3 | 2.000 | 83.89 | 0 | 78.94 | 0 | 21-133 | 6 | 0-25 | 3 |
| 2-Methylnaphthalene | 42.74 | 2.000 | 34.71 | 0 | 32.01 | 0 | 21-140 | 8 | 0-25 | 3 |
| 1-Methylnaphthalene | 59.00 | 2.000 | 47.61 | 0 | 45.00 | 0 | 20-140 | 6 | 0-25 | 3 |
| Acenaphthylene | ND | 2.000 | 1.706 | 85 | 1.549 | 77 | 33-145 | 10 | 0-25 | |
| Acenaphthene | 0.5300 | 2.000 | 2.109 | 79 | 1.970 | 72 | 49-121 | 7 | 0-25 | |
| Fluorene | ND | 2.000 | 1.860 | 93 | 1.701 | 85 | 59-121 | 9 | 0-25 | |
| Phenanthrene | ND | 2.000 | 1.624 | 81 | 1.459 | 73 | 54-120 | 11 | 0-25 | |
| Anthracene | ND | 2.000 | 1.548 | 77 | 1.417 | 71 | 27-133 | 9 | 0-25 | |
| Fluoranthene | ND | 2.000 | 1.666 | 83 | 1.524 | 76 | 26-137 | 9 | 0-25 | |
| Pyrene | ND | 2.000 | 1.581 | 79 | 1.473 | 74 | 18-168 | 7 | 0-25 | |
| Benzo (a) Anthracene | ND | 2.000 | 1.551 | 78 | 1.422 | 71 | 33-143 | 9 | 0-25 | |
| Chrysene | ND | 2.000 | 1.540 | 77 | 1.402 | 70 | 17-168 | 9 | 0-25 | |
| Benzo (k) Fluoranthene | ND | 2.000 | 1.447 | 72 | 1.319 | 66 | 24-159 | 9 | 0-25 | |
| Benzo (b) Fluoranthene | ND | 2.000 | 1.372 | 69 | 1.248 | 62 | 24-159 | 9 | 0-25 | |
| Benzo (a) Pyrene | ND | 2.000 | 1.331 | 67 | 1.216 | 61 | 17-163 | 9 | 0-25 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 2.000 | 1.320 | 66 | 1.166 | 58 | 10-171 | 12 | 0-25 | |
| Dibenz (a,h) Anthracene | ND | 2.000 | 1.422 | 71 | 1.275 | 64 | 10-219 | 11 | 0-25 | |
| Benzo (g,h,i) Perylene | ND | 2.000 | 1.524 | 76 | 1.373 | 69 | 10-227 | 10 | 0-25 | |



GC/MS / EPA 8260B



Quality Control - Spike/Spike Duplicate

Environmental Science International, Inc.

Date Received:

Work Order:

10/28/14

Work Order:

Kailua, HI 96734-2500

Preparation:

EPA 5030C

Method:

Project: Red Hill LTM 112066 Page 4 of 5

| | | | Matrix | | nstrument | Date Prepare | u Date Ana | lyzed | M2/M2D Ra | itch Number |
|-----------------------------|-------------------------------|-----------------------|-------------|------------|--------------|--------------|------------|-------|------------|-------------|
| ES114 | Sample | | Aqueou | s (| GC/MS OO | 10/28/14 | 10/29/14 | 05:51 | 1410285037 | 7 |
| ES114 | Matrix Spike | | Aqueou | s (| GC/MS OO | 10/28/14 | 10/29/14 | 07:40 | 141028S037 | 7 |
| ES114 | Matrix Spike | Duplicate | Aqueou | s (| 3C/MS OO | 10/28/14 | 10/29/14 | 08:07 | 1410285037 | 7 |
| <u>Parameter</u> | <u>Sample</u> <u>Conc.</u> | <u>Spike</u> Added | MS Conc. | MS %Rec | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Acetone | ND | 50.00 | 55.96 | 112 | 59.02 | 118 | 40-140 | 5 | 0-20 | |
| Benzene | ND | 50.00 | 48.83 | 98 | 47.07 | 94 | 80-120 | 4 | 0-20 | |
| Bromodichloromethane | ND | 50.00 | 48.14 | 96 | 47.21 | 94 | 75-120 | 2 | 0-20 | |
| Bromoform | ND | 50.00 | 46.72 | 93 | 46.10 | 92 | 70-130 | 1 | 0-20 | |
| Bromomethane | ND | 50.00 | 47.90 | 96 | 42.68 | 85 | 30-145 | 12 | 0-20 | |
| 2-Butanone | ND | 50.00 | 49.72 | 99 | 54.71 | 109 | 30-150 | 10 | 0-20 | |
| Carbon Tetrachloride | ND | 50.00 | 41.49 | 83 | 41.52 | 83 | 65-140 | 0 | 0-20 | |
| Chlorobenzene | ND | 50.00 | 49.02 | 98 | 47.13 | 94 | 80-120 | 4 | 0-20 | |
| Chloroethane | ND | 50.00 | 48.97 | 98 | 46.89 | 94 | 60-135 | 4 | 0-20 | |
| Chloroform | ND | 50.00 | 48.66 | 97 | 47.25 | 94 | 65-135 | 3 | 0-20 | |
| Chloromethane | ND | 50.00 | 42.58 | 85 | 42.98 | 86 | 40-125 | 1 | 0-20 | |
| Dibromochloromethane | ND | 50.00 | 49.94 | 100 | 48.76 | 98 | 60-135 | 2 | 0-20 | |
| 1,2-Dibromo-3-Chloropropane | ND | 50.00 | 42.13 | 84 | 42.15 | 84 | 50-130 | 0 | 0-20 | |
| 1,2-Dibromoethane | ND | 50.00 | 50.33 | 101 | 49.09 | 98 | 80-120 | 2 | 0-20 | |
| 1,2-Dichlorobenzene | ND | 50.00 | 47.77 | 96 | 46.40 | 93 | 70-120 | 3 | 0-20 | |
| 1,3-Dichlorobenzene | ND | 50.00 | 47.15 | 94 | 45.84 | 92 | 75-125 | 3 | 0-20 | |
| 1,4-Dichlorobenzene | ND | 50.00 | 46.75 | 94 | 45.19 | 90 | 75-125 | 3 | 0-20 | |
| 1,1-Dichloroethane | ND | 50.00 | 45.35 | 91 | 44.62 | 89 | 70-135 | 2 | 0-20 | |
| 1,2-Dichloroethane | ND | 50.00 | 54.76 | 110 | 52.87 | 106 | 70-130 | 4 | 0-20 | |
| 1,1-Dichloroethene | ND | 50.00 | 46.83 | 94 | 46.17 | 92 | 70-130 | 1 | 0-20 | |
| c-1,2-Dichloroethene | ND | 50.00 | 48.30 | 97 | 47.53 | 95 | 70-125 | 2 | 0-20 | |
| t-1,2-Dichloroethene | ND | 50.00 | 46.93 | 94 | 46.11 | 92 | 60-140 | 2 | 0-20 | |
| 1,2-Dichloropropane | ND | 50.00 | 47.05 | 94 | 46.05 | 92 | 75-125 | 2 | 0-20 | |
| c-1,3-Dichloropropene | ND | 50.00 | 43.96 | 88 | 43.95 | 88 | 70-130 | 0 | 0-20 | |
| t-1,3-Dichloropropene | ND | 50.00 | 37.58 | 75 | 38.14 | 76 | 55-140 | 1 | 0-20 | |
| Ethylbenzene | ND | 50.00 | 48.57 | 97 | 46.29 | 93 | 75-125 | 5 | 0-20 | |
| Methylene Chloride | ND | 50.00 | 46.74 | 93 | 46.27 | 93 | 55-140 | 1 | 0-20 | |
| 4-Methyl-2-Pentanone | ND | 50.00 | 45.66 | 91 | 46.54 | 93 | 60-135 | 2 | 0-20 | |
| Styrene | ND | 50.00 | 49.11 | 98 | 47.51 | 95 | 65-135 | 3 | 0-20 | |
| 1,1,1,2-Tetrachloroethane | ND | 50.00 | 46.10 | 92 | 45.50 | 91 | 80-130 | 1 | 0-20 | |
| 1,1,2,2-Tetrachloroethane | ND | 50.00 | 48.52 | 97 | 47.00 | 94 | 65-130 | 3 | 0-20 | |
| Tetrachloroethene | ND | 50.00 | 53.01 | 106 | 50.86 | 102 | 45-150 | 4 | 0-20 | |
| Toluene | ND | 50.00 | 48.79 | 98 | 47.39 | 95 | 75-120 | 3 | 0-20 | |
| 1,2,4-Trichlorobenzene | ND | 50.00 | 47.10 | 94 | 46.48 | 93 | 65-135 | 1 | 0-20 | |
| 1,1,1-Trichloroethane | ND | 50.00 | 43.81 | 88 | 43.80 | 88 | 65-130 | 0 | 0-20 | |



Quality Control - Spike/Spike Duplicate

Environmental Science International, Inc.

Date Received:

Work Order:

14-10-2131

Kailua, HI 96734-2500

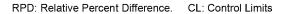
Preparation:

Method:

GC/MS / EPA 8260B

Project: Red Hill LTM 112066 Page 5 of 5

| <u>Parameter</u> | <u>Sample</u> <u>Conc.</u> | <u>Spike</u> Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|-----------------------------|-------------------------------|-----------------------|-------------|-------------|--------------|--------------|----------|-----|--------|------------|
| Hexachloro-1,3-Butadiene | ND | 50.00 | 47.57 | 95 | 45.21 | 90 | 50-140 | 5 | 0-20 | |
| 1,1,2-Trichloroethane | ND | 50.00 | 47.82 | 96 | 46.98 | 94 | 75-125 | 2 | 0-20 | |
| Trichloroethene | ND | 50.00 | 47.24 | 94 | 44.84 | 90 | 70-125 | 5 | 0-20 | |
| 1,2,3-Trichloropropane | ND | 50.00 | 50.56 | 101 | 49.71 | 99 | 75-125 | 2 | 0-20 | |
| Vinyl Chloride | ND | 50.00 | 51.75 | 103 | 51.09 | 102 | 50-145 | 1 | 0-20 | |
| p/m-Xylene | ND | 100.0 | 96.07 | 96 | 92.12 | 92 | 75-130 | 4 | 0-20 | |
| o-Xylene | ND | 50.00 | 49.17 | 98 | 47.56 | 95 | 80-120 | 3 | 0-20 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 50.00 | 40.87 | 82 | 42.14 | 84 | 65-125 | 3 | 0-20 | |



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Project: Red Hill LTM 112066

Quality Control - PDS

Environmental Science International, Inc.

Date Received:

Work Order:

10/28/14

Work Order:

Kailua, HI 96734-2500

Preparation:

Method:

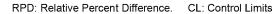
EPA 3005A Filt.

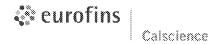
Method:

Page 1 of 1

| Quality Control Sample ID | Туре | N | latrix | Instrument | Date Prepared Date | e Analyzed PDS Num | |
|---------------------------|--------|--------------|-------------|------------|---------------------|-----------------------|------------|
| ES114 | Sample | А | queous | ICP/MS 04 | 10/29/14 00:00 10/3 | 10/14 23:32 1411 | 029S03B |
| ES114 | PDS | А | queous | ICP/MS 04 | 10/29/14 00:00 10/3 | 0/14 23:22 141 | 029S03B |
| <u>Parameter</u> | | Sample Conc. | Spike Added | PDS Conc. | PDS %Rec. | %Rec. CL | Qualifiers |
| Lead | | ND | 100.0 | 106.6 | 107 | 75-125 | |

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Quality Control - LCS/LCSD

Environmental Science International, Inc.

Date Received:

Work Order:

14-10-2131

Kailua, HI 96734-2500

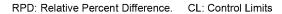
Preparation:

Method:

EPA 8015B (M)

Project: Red Hill LTM 112066 Page 1 of 5

| Quality Control Sample ID | Туре | Mat | rix | Instrument | Date Pre | pared Date | Analyzed | LCS/LCSD B | atch Number |
|---------------------------|-------------|-----------|--------------|------------|---------------|------------|------------|------------|-------------------|
| 099-15-516-206 | LCS | Aqı | ieous | GC 45 | 10/30/14 | 10/3 | 1/14 07:20 | 141030B06 | |
| 099-15-516-206 | LCSD | Aqu | ieous | GC 45 | 10/30/14 | 10/3 | 1/14 07:37 | 141030B06 | |
| Parameter | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | <u>Qualifiers</u> |
| TPH as Diesel | 2000 | 2296 | 115 | 2219 | 111 | 60-132 | 3 | 0-11 | |





Quality Control - LCS

Environmental Science International, Inc.

Date Received:

Work Order:

10/28/14

Work Order:

14-10-2131

Kailua, HI 96734-2500

Preparation:

Method:

EPA 3005A Filt.

Method:

Project: Red Hill LTM 112066 Page 2 of 5

| Quality Control Sample ID | Туре | Matrix | Instrument Da | | ate Analyzed LCS Bat | |
|---------------------------|------|-------------|-----------------|-----------|-----------------------|-------------------|
| 099-14-497-101 | LCS | Aqueous | ICP/MS 04 10 | /29/14 10 | 0/30/14 23:02 141029L | .03D |
| <u>Parameter</u> | | Spike Added | Conc. Recovered | LCS %Rec. | <u>%Rec. CL</u> | <u>Qualifiers</u> |
| Lead | | 100.0 | 106.3 | 106 | 80-120 | |





Quality Control - LCS

Environmental Science International, Inc.

354 Uluniu Street, Suite 304 Kailua, HI 96734-2500

Project: Red Hill LTM 112066

Date Received:

10/28/14 14-10-2131

Work Order: Preparation:

EPA 3510C EPA 8270C SIM PAHs

Method:

Page 3 of 5

| Quality Control Sample ID | Туре | Matrix | Instrument [| Date Prepared Date | e Analyzed LCS Ba | atch Number |
|---------------------------|------|-------------|----------------|--------------------|----------------------|-------------|
| 099-15-148-67 | LCS | Aqueous | GC/MS AAA 1 | 10/30/14 10/3 | 31/14 17:40 141030 | L08 |
| <u>Parameter</u> | | Spike Added | Conc. Recovere | d LCS %Rec. | %Rec. CL | Qualifiers |
| Naphthalene | | 2.000 | 1.394 | 70 | 21-133 | |
| 2-Methylnaphthalene | | 2.000 | 1.382 | 69 | 21-140 | |
| 1-Methylnaphthalene | | 2.000 | 1.409 | 70 | 20-140 | |
| Acenaphthylene | | 2.000 | 1.394 | 70 | 33-145 | |
| Acenaphthene | | 2.000 | 1.451 | 73 | 55-121 | |
| Fluorene | | 2.000 | 1.539 | 77 | 59-121 | |
| Phenanthrene | | 2.000 | 1.447 | 72 | 54-120 | |
| Anthracene | | 2.000 | 1.308 | 65 | 27-133 | |
| Fluoranthene | | 2.000 | 1.538 | 77 | 26-137 | |
| Pyrene | | 2.000 | 1.480 | 74 | 45-129 | |
| Benzo (a) Anthracene | | 2.000 | 1.398 | 70 | 33-143 | |
| Chrysene | | 2.000 | 1.425 | 71 | 17-168 | |
| Benzo (k) Fluoranthene | | 2.000 | 1.412 | 71 | 24-159 | |
| Benzo (b) Fluoranthene | | 2.000 | 1.322 | 66 | 24-159 | |
| Benzo (a) Pyrene | | 2.000 | 1.150 | 58 | 17-163 | |
| Indeno (1,2,3-c,d) Pyrene | | 2.000 | 1.251 | 63 | 25-175 | |
| Dibenz (a,h) Anthracene | | 2.000 | 1.391 | 70 | 25-175 | |
| Benzo (g,h,i) Perylene | | 2.000 | 1.514 | 76 | 25-157 | |



10/28/14

14-10-2131

EPA 5030C

GC/MS / EPA 8260B



Quality Control - LCS/LCSD

Environmental Science International, Inc.

Date Received:

Work Order:

Kailua, HI 96734-2500 Preparation:

Project: Red Hill LTM 112066 Page 4 of 5

Method:

| Quality Control Sample ID | Туре | Mat | rix | Instrument | Date Pre | epared [| Date Analyzed | LCS/LCSD B | atch Number |
|-----------------------------|-------------|-----------|---------------------|------------|---------------|----------|----------------|------------|-------------|
| 099-13-057-71 | LCS | Aqı | ieous | GC/MS OO | 10/28/14 | | 10/29/14 03:34 | 141028L046 | |
| 099-13-057-71 | LCSD | Aqı | ieous | GC/MS OO | 10/28/14 | | 10/29/14 04:01 | 141028L046 | |
| <u>Parameter</u> | Spike Added | LCS Conc. | <u>LCS</u> %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. | CL RPD | RPD CL | Qualifiers |
| Acetone | 50.00 | 44.78 | 90 | N/A | N/A | 40-140 | N/A | 0-20 | |
| Benzene | 50.00 | 51.34 | 103 | N/A | N/A | 80-120 | N/A | 0-20 | |
| Bromodichloromethane | 50.00 | 51.61 | 103 | N/A | N/A | 75-120 | N/A | 0-20 | |
| Bromoform | 50.00 | 49.35 | 99 | N/A | N/A | 70-130 | N/A | 0-20 | |
| Bromomethane | 50.00 | 44.25 | 88 | N/A | N/A | 30-145 | N/A | 0-20 | |
| 2-Butanone | 50.00 | 47.46 | 95 | N/A | N/A | 30-150 | N/A | 0-20 | |
| Carbon Tetrachloride | 50.00 | 45.03 | 90 | N/A | N/A | 65-140 | N/A | 0-20 | |
| Chlorobenzene | 50.00 | 50.81 | 102 | N/A | N/A | 80-120 | N/A | 0-20 | |
| Chloroethane | 50.00 | 52.54 | 105 | N/A | N/A | 60-135 | N/A | 0-20 | |
| Chloroform | 50.00 | 51.69 | 103 | N/A | N/A | 65-135 | N/A | 0-20 | |
| Chloromethane | 50.00 | 44.16 | 88 | N/A | N/A | 40-125 | N/A | 0-20 | |
| Dibromochloromethane | 50.00 | 52.37 | 105 | N/A | N/A | 60-135 | N/A | 0-20 | |
| 1,2-Dibromo-3-Chloropropane | 50.00 | 41.53 | 83 | N/A | N/A | 50-130 | N/A | 0-20 | |
| 1,2-Dibromoethane | 50.00 | 52.74 | 105 | N/A | N/A | 80-120 | N/A | 0-20 | |
| 1,2-Dichlorobenzene | 50.00 | 49.45 | 99 | N/A | N/A | 70-120 | N/A | 0-20 | |
| 1,3-Dichlorobenzene | 50.00 | 48.61 | 97 | N/A | N/A | 75-125 | N/A | 0-20 | |
| 1,4-Dichlorobenzene | 50.00 | 48.44 | 97 | N/A | N/A | 75-125 | N/A | 0-20 | |
| 1,1-Dichloroethane | 50.00 | 48.98 | 98 | N/A | N/A | 70-135 | N/A | 0-20 | |
| 1,2-Dichloroethane | 50.00 | 58.50 | 117 | N/A | N/A | 70-130 | N/A | 0-20 | |
| 1,1-Dichloroethene | 50.00 | 50.88 | 102 | N/A | N/A | 70-130 | N/A | 0-20 | |
| c-1,2-Dichloroethene | 50.00 | 51.67 | 103 | N/A | N/A | 70-125 | N/A | 0-20 | |
| t-1,2-Dichloroethene | 50.00 | 50.30 | 101 | N/A | N/A | 60-140 | N/A | 0-20 | |
| 1,2-Dichloropropane | 50.00 | 50.03 | 100 | N/A | N/A | 75-125 | N/A | 0-20 | |
| c-1,3-Dichloropropene | 50.00 | 48.68 | 97 | N/A | N/A | 70-130 | N/A | 0-20 | |
| t-1,3-Dichloropropene | 50.00 | 43.01 | 86 | N/A | N/A | 55-140 | N/A | 0-20 | |
| Ethylbenzene | 50.00 | 50.41 | 101 | N/A | N/A | 75-125 | N/A | 0-20 | |
| Methylene Chloride | 50.00 | 50.23 | 100 | N/A | N/A | 55-140 | N/A | 0-20 | |
| 4-Methyl-2-Pentanone | 50.00 | 49.12 | 98 | N/A | N/A | 60-135 | N/A | 0-20 | |
| Styrene | 50.00 | 50.88 | 102 | N/A | N/A | 65-135 | N/A | 0-20 | |
| 1,1,1,2-Tetrachloroethane | 50.00 | 48.61 | 97 | N/A | N/A | 80-130 | N/A | 0-20 | |
| 1,1,2,2-Tetrachloroethane | 50.00 | 49.80 | 100 | N/A | N/A | 65-130 | N/A | 0-20 | |
| Tetrachloroethene | 50.00 | 52.93 | 106 | N/A | N/A | 45-150 | N/A | 0-20 | |
| Toluene | 50.00 | 51.72 | 103 | N/A | N/A | 75-120 | N/A | 0-20 | |
| 1,2,4-Trichlorobenzene | 50.00 | 48.11 | 96 | N/A | N/A | 65-135 | | 0-20 | |
| 1,1,1-Trichloroethane | 50.00 | 47.64 | 95 | N/A | N/A | 65-130 | | 0-20 | |
| Hexachloro-1,3-Butadiene | 50.00 | 48.39 | 97 | N/A | N/A | 50-140 | | 0-20 | |
| | | | | | | | | | |



Quality Control - LCS/LCSD

Environmental Science International, Inc.

Date Received:

Work Order:

14-10-2131

Kailua, HI 96734-2500

Preparation:

Method:

Date Received:

10/28/14

Preparation:

EPA 5030C

Project: Red Hill LTM 112066 Page 5 of 5

| <u>Parameter</u> | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|-----------------------------|-------------|-----------|--------------|------------|---------------|----------|-----|--------|------------|
| 1,1,2-Trichloroethane | 50.00 | 50.32 | 101 | N/A | N/A | 75-125 | N/A | 0-20 | |
| Trichloroethene | 50.00 | 49.50 | 99 | N/A | N/A | 70-125 | N/A | 0-20 | |
| 1,2,3-Trichloropropane | 50.00 | 52.64 | 105 | N/A | N/A | 75-125 | N/A | 0-20 | |
| Vinyl Chloride | 50.00 | 54.73 | 109 | N/A | N/A | 50-145 | N/A | 0-20 | |
| p/m-Xylene | 100.0 | 100.3 | 100 | N/A | N/A | 75-130 | N/A | 0-20 | |
| o-Xylene | 50.00 | 51.10 | 102 | N/A | N/A | 80-120 | N/A | 0-20 | |
| Methyl-t-Butyl Ether (MTBE) | 50.00 | 45.38 | 91 | N/A | N/A | 65-125 | N/A | 0-20 | |
| Gasoline Range Organics | 1000 | 1077 | 108 | 1032 | 103 | 80-120 | 4 | 0-20 | |

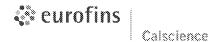




Sample Analysis Summary Report

| Work Order: 14-10-2131 | Work Order: 14-10-2131 | | | | | | |
|------------------------|------------------------|------------|------------|---------------------|--|--|--|
| <u>Method</u> | Extraction | Chemist ID | Instrument | Analytical Location | | | |
| EPA 6020 | EPA 3005A Filt. | 598 | ICP/MS 04 | 1 | | | |
| EPA 8015B (M) | EPA 3510C | 682 | GC 45 | 1 | | | |
| EPA 8270C SIM PAHs | EPA 3510C | 923 | GC/MS AAA | 1 | | | |
| GC/MS / EPA 8260B | EPA 5030C | 849 | GC/MS OO | 2 | | | |

Location 1: 7440 Lincoln Way, Garden Grove, CA 92841 Location 2: 7445 Lampson Avenue, Garden Grove, CA 92841



Glossary of Terms and Qualifiers

Work Order: 14-10-2131 Page 1 of 1

| Qualifiers | <u>Definition</u> |
|------------|--|
| * | See applicable analysis comment. |
| < | Less than the indicated value. |
| > | Greater than the indicated value. |
| 1 | Surrogate compound recovery was out of control due to a required sample dilution. Therefore, the sample data was reported without further clarification. |
| 2 | Surrogate compound recovery was out of control due to matrix interference. The associated method blank surrogate spike compound was in control and, therefore, the sample data was reported without further clarification. |
| 3 | Recovery of the Matrix Spike (MS) or Matrix Spike Duplicate (MSD) compound was out of control due to suspected matrix interference. The associated LCS recovery was in control. |
| 4 | The MS/MSD RPD was out of control due to suspected matrix interference. |
| 5 | The PDS/PDSD or PES/PESD associated with this batch of samples was out of control due to suspected matrix interference. |
| 6 | Surrogate recovery below the acceptance limit. |
| 7 | Surrogate recovery above the acceptance limit. |
| В | Analyte was present in the associated method blank. |
| BU | Sample analyzed after holding time expired. |
| BV | Sample received after holding time expired. |
| DL | The Detection Limit (DL) is the smallest analyte concentration that can be demonstrated to be different from zero or a blank concentration at the 99% level of confidence. |
| Е | Concentration exceeds the calibration range. |
| ET | Sample was extracted past end of recommended max. holding time. |
| HD | The chromatographic pattern was inconsistent with the profile of the reference fuel standard. |
| ICH | Initial calibration verification recovery is above the control limit for this analyte. |
| ICJ | Initial calibration verification recovery is below the control limit for this analyte. |
| IH | Calibration verification recovery is above the control limit for this analyte. |
| IJ | Calibration verification recovery is below the control limit for this analyte. |
| J | Analyte was detected at a concentration below the LOQ and above the DL. Reported value is estimated. |
| JA | Analyte positively identified but quantitation is an estimate. |
| LOD | The Limit of Detection (LOD) is the smallest amount or concentration of a substance that must be present in a sample in order to be detected at 99% confidence level. |
| LOQ | The Limit of Quantitation (LOQ) is the lowest concentration of a substance that produces a quantitative result within specified limits of precision and bias. |
| Q | Spike recovery and RPD control limits do not apply resulting from the parameter concentration in the sample exceeding the spike concentration by a factor of four or greater. |
| SG | The sample extract was subjected to Silica Gel treatment prior to analysis. |
| U | Undetected at Detection Limit (DL) and is reported as less than the Limit of Detection (LOD). |
| Χ | % Recovery and/or RPD out-of-range. |
| Z | Analyte presence was not confirmed by second column or GC/MS analysis. |
| | Solid - Unless otherwise indicated, solid sample data is reported on a wet weight basis, not corrected for % moisture. All QC results are |

reported on a wet weight basis.

Any parameter identified in 40CFR Part 136.3 Table II that is designated as "analyze immediately" with a holding time of <= 15 minutes (40CFR-136.3 Table II, footnote 4), is considered a "field" test and the reported results will be qualified as being received outside of the stated holding time unless received at the laboratory within 15 minutes of the collection time.

A calculated total result (Example: Total Pesticides) is the summation of each component concentration and/or, if "J" flags are reported, estimated concentration. Component concentrations showing not detected (ND) are summed into the calculated total result as zero concentrations.

| Calscience Environmental Laboratories 7440 Lincoln Way, Garden Grove, CA 92841-1427 • (714) 895-5494 | ence Environmental Laboratorie 7440 Lincoln Way, Garden Grove, CA 92841-1427 (714) 895-5494 | Laboral -1427 • (714) | COMPOS 395-5494 | | Same and the same | 0#/LAI | WO#/LAB USE ONLY | NLY | | | Date | 3 | | J E | 32 | CUSTODY RECORD | | |
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| Other locations: Col | Other locations: Concord, San Luis Obispo, Houston, and Corpus Christi For courier service / sample drop off information, contact <u>sales@calscience.com</u> or call us. | ouston, and Colf information, or call us. | rpus Christi | | * | | | | | | Page | (1) | | | 10 J | or the rest of the second seco | | 1 1 |
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SHIP DATE: 270CT14 ACTWGT: 56.1 LB CAD: /POS1525 DIMS: 26x14x13 IN

BILL RECIPIENT

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GARDEN GROVE CA 92841

FedEx Express

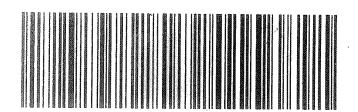


1 of 2 TRK# 0200 8045 5791 7364 ## MASTER ##

WZ APVA

TUE - 28 OCT AA STANDARD OVERNIGHT

> 92841 CA-US SNA



ORIGÍN ID: HÑLA

SHIP DATE: 270CT14 ACTWGT: 48,4 LB CAD: /POS1525 DIMS: 26x14x13 IN

BILL RECIPIENT

UNITED STATES US

SAMPLE CONTROL CALSCIENCE 7440 LINCOLN WAY

GARDEN GROVE CA 92841

REF:

FedEx Express

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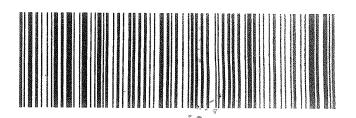
2 of 2 MPS# 7801 5009 0734 Mstr# 8045 5791 7364

TUE - 28 OCT STANDARD OVERNIGHT

0200

APVA

92841 CA-US SNA





Calscience

WORK ORDER #: 14-10- 2 1 3 1

SAMPLE RECEIPT FORM

| Cooler | | of | 2_ |
|----------|-----|----|----|
| 10 / | n C | 11 | Λ |

| CLIENT: | SE | | | DATE: _ | 10/28 | <u>/ 14</u> |
|--|---|--|---|------------------------|-------------------|---|
| TEMPERATURE: | Thermometer | ID: SC2 (Criteria: 0.0 °C | – 6.0 °C, not froz | en except se | diment/tissu | e) |
| 9 | | - 0.2°C (CF) = | | Blank | ☐ Sampl | |
| | | criteria (PM/APM contact | | | ŕ | |
| | | criteria but received on ic | | dav of sampl | ina. | |
| | | ture, placed on ice fo | | | 9. | |
| Ambient Temperate | - | | r trunsport by c | ounon | Checked b | v: 15 |
| Ambient remperati | uic. Li All | ☐ Filter | | | | |
| CUSTODY SEALS | INTACT: | | | | | |
| © Cooler □ | | □ No (Not Intact) | ☐ Not Presen | t □ N/A | Checked b | y: <u>LS</u> |
| □Sample □ | | ☐ No (Not Intact) | □ Not Presen | t | Checked b | y: <u>876</u> |
| | | | | | | |
| SAMPLE CONDIT | | | | Yes | No | N/A |
| | | nt(s) received with san | | | | |
| | | ete | | | | |
| | | # of containers logged in ba | | S. | | |
| ☐ No analysis reques | | , | me relinquished. | | | |
| 8 | | D | | | | |
| 1 | ` , | ent with COC | | * | . П | |
| | | od condition | | | | |
| | | olume for analyses req | | | | |
| | * | ime | | | | |
| | | hin 15-minute holding t | | | [] | |
| ' | | ssolved Sulfides | | _ | | |
| , , | | C or sample container. | , | | | |
| Unpreserved vial | | | | | $\overline{}$ | |
| | | of headspace | | | | |
| CONTAINER TYP | E: | n | | | | |
| Solid: □4ozCGJ | □8ozCGJ □ | I16ozCGJ □Sleeve (|) □EnCor | es [®] □Terra | aCores® □_ | |
| Aqueous: 🗹 VOA | ÍVOAK □VOA | A na₂ □125AGB □125 <i>i</i> | AGB h □125AGB | p ZÍAGB | □1AGB na ₂ | □1AGB s |
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| Air: □Tedlar® □Ca Container: C: Clear A: Ar | anister Other: mber P : Plastic G : | Trip Blank Glass J: Jar B: Bottle Z: Ziplor a: NaOH p: H ₃ PO ₄ s: H ₂ SO ₄ u: U | K Lot#: <u>1404037</u> c/Resealable Bag E: | Labeled Envelope | Reviewed by | : _07 |



Calscience

WORK ORDER #: 14-10- 2 1 5

SAMPLE RECEIPT FORM

Cooler 2 of 2

| CLIENT: EST | DATE:_ | 10/28 | /14 |
|---|--|-----------------------------|----------------|
| TEMPERATURE: Thermometer ID: SC2 (Criteria: 0.0 °C – 6.0 °C, not t | rozen except se | diment/tissu | e) |
| Temperature °C - 0.2 °C (CF) = 4 °C | AND THE RESERVE TO THE PERSON NAMED TO THE PER | ☐ Sample | |
| ☐ Sample(s) outside temperature criteria (PM/APM contacted by: | | • | |
| ☐ Sample(s) outside temperature criteria but received on ice/chilled on sa | | ina | |
| ☐ Received at ambient temperature, placed on ice for transport b | | 9. | |
| Ambient Temperature: | y Courier. | Checked by | v. 15 |
| Ambient reimperature. | | Officer D | y· |
| CUSTODY SEALS INTACT: | | | |
| Cooler □ □ No (Not Intact) □ Not Pres | sent □ N/A | Checked by | y: <u>LS</u> |
| Sample | sent | Checked by | 1: _836_ |
| CAMBLE CONDITION. | V | N. C. | NI/Λ |
| SAMPLE CONDITION: | Yes | No | N/A |
| Chain-Of-Custody (COC) document(s) received with samples | | | |
| COC document(s) received complete | | | |
| □ Collection date/time, matrix, and/or # of containers logged in based on sample land to the containers logged in based on the containers logged in b | | | |
| Sampler's name indicated on COC | | | |
| Sample container label(s) consistent with COC | | | |
| Sample container(s) intact and good condition | | | |
| Proper containers and sufficient volume for analyses requested | | П | |
| Analyses received within holding time | / | | П |
| Agueous samples received within 15-minute holding time | | | - |
| ☐ pH ☐ Residual Chlorine ☐ Dissolved Sulfides ☐ Dissolved Oxygen | 🗆 | | |
| Proper preservation noted on COC or sample container | _ | | |
| ☐ Unpreserved vials received for Volatiles analysis | | | |
| Volatile analysis container(s) free of headspace | | | |
| Tedlar bag(s) free of condensation | | | |
| CONTAINER TYPE: | | | |
| Solid: 4ozCGJ 8ozCGJ 16ozCGJ Sleeve () En | | | |
| Aqueous: □VOA □VOAh □VOAna₂ □125AGB □125AGBh □125A | GB p 21AGB [| ∃1AGB na ₂ [|]1AGB s |
| □500AGB ☑500AGJ □500AGJs □250AGB □250CGB □250C | GBs □1PB | □1PB na □ | 500PB |
| □250PB ଢ250PBn □125PB □125PB znna □100PJ □100PJ na ₂ | | | |
| Air: □Tedlar [®] □Canister Other : □ Trip Blank Lot#: | | | |
| Container: C: Clear A: Amber P: Plastic G: Glass J: Jar B: Bottle Z: Ziploc/Resealable Bag | • | Reviewed by: Scanned by: | f # [|

RAW DATA SHEET FOR METHOD: EPA 8015B (M)

WORK ORDER: INSTRUMENT:

14-10-2131

GC 45

EPA 3510C EXTRACTION: D/T EXTRACTED:

2014-10-30 00:00

ANALYZED BY:

682 D/T ANALYZED:

2014-10-31 09:41

REVIEWED BY: 628

2014-11-04 17:13 D/T REVIEWED:

DATA FILE:

W:\GC 45\GC 45 DATA\2014\141030\14103063.D\14103063

2

CLIENT SAMPLE NUMBER: ES113

LCS/MB BATCH: MS/MSD BATCH:

141030B06

SAMPLE VOLUME / WEIGHT:

DEFAULT: 500.00 ml / ACTUAL: 500.00 ml

DEFAULT: 5.00 ml / ACTUAL: 2.50 ml

LOD

UNITS:

ug/L

FINAL VOLUME / WEIGHT: ADJUSTMENT RATIO TO PF:

0.50

COMMENT:

Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are

qualified with a "J" flag. TPH as Diesel is quantified in the carbon range C10-C28.

COMPOUND

INI. CONC

DF

CONC

DL 11

LOQ

QUAL

TPH as Diesel

1.00 23400

117

12

25

b

Data File Name : W:\GC_45\GC 45 DATA\2014\141030\14103063.D

Page Number : 1

Operator : 682 Vial Number : Vial 63

Instrument : GC 45 Injection Number : 1
Sample Name : 14-10-2131-2 Sequence Line : 63

Instrument Method: C:\CHEM32\1\METHODS\ ->

Acquired on : 31 Oct 14 09:41 am

Report Created on: 04 Nov 14 02:48 pm Analysis Method : 8015B.MTH

Software Revision: Rev. B.03.02 [341] Copyright © Agilent Technologies

Sig. 1 in W:\GC 45\GC 45 DATA\2014\141030\ ->

| Pk | Ret Time | Area | Height | Peak | | Response % |
|----|----------|--------|--------|------|-------|------------|
| 1 | 2.024 | 2.51 | | VV | 0.035 | 0.239 |
| 2 | 2.092 | 1.31 | | VV | 0.048 | 0.125 |
| 3 | 2.223 | 7.06 | | VV | 0.039 | 0.671 |
| 4 | 2.285 | 1.69 | | VV | 0.022 | 0.161 |
| 5 | 2.334 | 3.26 | | VV | 0.034 | 0.309 |
| 6 | 2.463 | 10.00 | 3 | VV | 0.048 | 0.950 |
| 7 | 2.511 | 3.68 | 2 | VV | 0.028 | 0.349 |
| 8 | 2.638 | 13.63 | 3 | vv | 0.070 | 1.294 |
| 9 | 2.753 | 24.19 | 5 | VV | 0.060 | 2.297 |
| 10 | 2.883 | 24.76 | 7 | VV | 0.045 | 2.351 |
| 11 | 2.943 | 15.15 | 5 | VV | 0.044 | 1.439 |
| 12 | 3.012 | 16.40 | 4 | VV | 0.062 | 1.558 |
| 13 | 3.094 | 8.03 | 4 | VV | 0.032 | 0.762 |
| 14 | 3.168 | 19.46 | 4 | VV | 0.056 | 1.848 |
| 15 | 3.306 | 37.68 | 4 | VV | 0.109 | 3.578 |
| 16 | 3.438 | 14.10 | 4 | VV | 0.050 | 1.339 |
| 17 | 3.508 | 16.84 | 4 | VV | 0.059 | 1.599 |
| 18 | 3.559 | 10.28 | 4 | VV | 0.041 | 0.976 |
| 19 | 3.602 | 13.03 | 3 | VV | 0.048 | 1.238 |
| 20 | 3.715 | 25.80 | 4 | VV | 0.089 | 2.450 |
| 21 | 3.849 | 7.80 | 3 | VV | 0.040 | 0.741 |
| 22 | 3.911 | 9.29 | 2 | VV | 0.067 | 0.882 |
| 23 | 4.017 | 11.82 | | VV | 0.052 | 1.123 |
| 24 | 4.171 | 1.93 | | VV | 0.034 | 0.183 |
| 25 | 4.238 | 5.50 | | VV | 0.075 | 0.523 |
| 26 | 4.399 | 7.57 | | VV | 0.119 | 0.718 |
| 27 | 4.593 | 1.54 | | VV | 0.046 | 0.146 |
| 28 | 4.690 | 5.13 | | VV | 0.090 | 0.487 |
| 29 | 4.848 | 2.51 | | VV | 0.046 | |
| 30 | 4.909 | 1.46 | | VV | 0.036 | |
| 31 | 4.967 | 1.29 | | VV | 0.034 | 0.122 |
| 32 | 5.069 | 3.16 | | VV | 0.044 | 0.300 |
| 33 | 5.220 | 2.72 | | VV | 0.044 | 0.258 |
| 34 | 5.278 | 1.24 | | VV | 0.044 | 0.118 |
| 35 | 5.373 | 721.21 | 503 | VV | 0.024 | 68.488 |

Total area =

Data File Name : W:\GC_45\GC 45 DATA\2014\141030\14103063.D

Page Number : 2

Operator : 682 Vial Number : Vial 63

Instrument : GC 45 Injection Number : 1 Sample Name : 14-10-2131-2 Sequence Line : 63

Instrument Method: C:\CHEM32\1\METHODS\ -:

Acquired on : 31 Oct 14 09:41 am

Report Created on: 04 Nov 14 02:48 pm Analysis Method : 8015B.MTH

Software Revision: Rev. B.03.02 [341] Copyright © Agilent Technologies

1053.05

Data File Name : W:\GC 45\GC 45 DATA\2014\141030\14103063.D

Page Number : 3

Operator : 682 Vial Number : Vial 63

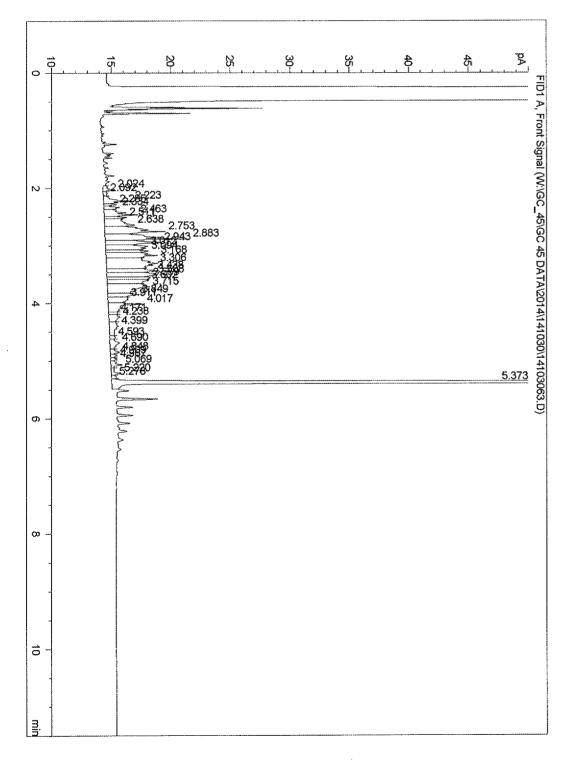
Instrument : GC 45 Injection Number : 1 Sample Name : 14-10-2131-2 Sequence Line : 63

Instrument Method: C:\CHEM32\1\METHODS\ ->

Acquired on : 31 Oct 14 09:41 am

Report Created on: 04 Nov 14 02:48 pm Analysis Method : 8015B.MTH

Software Revision: Rev. B.03.02 [341] Copyright © Agilent Technologies



RAW DATA SHEET FOR METHOD: EPA 8015B (M)

WORK ORDER:

14-10-2131

INSTRUMENT: GC 45

EPA 3510C

EXTRACTION: D/T EXTRACTED:

ANALYZED BY:

682 2014-10-31 09:59 D/T ANALYZED:

REVIEWED BY: 628

2014-10-30 00:00

D/T REVIEWED:

2014-11-04 17:13

DATA FILE:

W:\GC 45\GC 45 DATA\2014\141030\14103064.D\14103064

3

UNITS:

CLIENT SAMPLE NUMBER: ES114

LCS/MB BATCH: MS/MSD BATCH: 141030B06 141030S06

ug/L

SAMPLE VOLUME / WEIGHT:

DEFAULT: 500.00 ml / ACTUAL: 500.00 ml

DEFAULT: 5.00 ml / ACTUAL: 2.50 ml FINAL VOLUME / WEIGHT:

ADJUSTMENT RATIO TO PF: 0.50

COMMENT:

Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are

qualified with a "J" flag. TPH as Diesel is quantified in the carbon range C10-C28.

COMPOUND TPH as Diesel INI. CONC DF

409000

1.00

CONC 2050

LOD DL 11 12

LOQ

25

QUAL b

Area Percent Report

Data File Name : W:\GC_45\GC 45 DATA\2014\141030\14103064.D

Vial Number : Vial 64

Page Number : 1
Operator : 682
Instrument : GC Injection Number: 1 : GC 45 Sequence Line : 64 Sample Name : 14-10-2131-3

Instrument Method: C:\CHEM32\1\METHODS\

Acquired on : 31 Oct 14 09:59 am

Report Created on: 04 Nov 14 02:49 pm Analysis Method : 8015B.MTH

Software Revision: Rev. B.03.02 [341] Copyright © Agilent Technologies

Sig. 1 in W:\GC_45\GC 45 DATA\2014\141030\ ->

| Pk | Ret Time | Area | Height | | Width | Response % |
|----|----------|--------|--------|----|-------|------------|
| | | | | | | |
| 1 | 1.947 | 36.87 | | VV | 0.025 | 0.554 |
| 2 | 2.028 | 166.85 | 76 | | 0.032 | 2.505 |
| 3 | 2.173 | 220.62 | 128 | | 0.027 | 3.313 |
| 4 | 2.258 | 188.02 | 81 | VV | 0.032 | 2.823 |
| 5 | 2.353 | 227.05 | 110 | | 0.029 | 3.409 |
| б | 2.457 | 789.99 | 535 | | 0.022 | 11.862 |
| 7 | 2.544 | 191.88 | 47 | VV | 0.053 | 2,881 |
| 8 | 2.600 | 106.50 | 52 | VV | 0.030 | 1.599 |
| 9 | 2.642 | 116.77 | 54 | VV | 0.032 | 1.753 |
| 10 | 2.729 | 439.52 | 208 | VV | 0.030 | 6.599 |
| 11 | 2.769 | 366.68 | 248 | VV | 0.022 | 5.506 |
| 12 | 2.853 | 424.12 | 97 | VV | 0.056 | 6.368 |
| 13 | 2.948 | 289.07 | 81 | VV | 0.050 | 4.340 |
| 14 | 3.014 | 234.69 | 134 | VV | 0.025 | 3.524 |
| 15 | 3.058 | 266.23 | 81 | VV | 0.043 | 3.998 |
| 16 | 3.146 | 250.38 | 51 | VV | 0.064 | 3.759 |
| 17 | 3.298 | 368.20 | 51 | VV | 0.087 | 5.529 |
| 18 | 3.431 | 256.38 | 41 | VV | 0.083 | 3.850 |
| 19 | 3.498 | 130.37 | 37 | VV | 0.045 | 1.957 |
| 20 | 3.558 | 234.19 | 40 | VV | 0.074 | 3.516 |
| 21 | 3.687 | 134.10 | 32 | VV | 0.053 | 2.013 |
| 22 | 3.773 | 105.43 | 22 | VV | 0.060 | 1.583 |
| 23 | 3.853 | 46.10 | 18 | VV | 0.035 | 0.692 |
| 24 | 3.899 | 37.48 | 16 | VV | 0.032 | 0.563 |
| 25 | 3.938 | 58.23 | 19 | VV | 0.042 | 0.874 |
| 26 | 4.010 | 116.07 | 14 | VV | 0.102 | 1.743 |
| 27 | 4.261 | 19.12 | 4 | VV | 0.056 | 0.287 |
| 28 | 4.348 | 45.09 | 4 | VV | 0.154 | 0.677 |
| 29 | 4.681 | 20.80 | 3 | VV | 0.098 | 0.312 |
| 30 | 4.832 | 7.60 | 2 | VV | 0.050 | 0.114 |
| 31 | 4.891 | 5.24 | 2 | VV | 0.043 | 0.079 |
| 32 | 4.950 | 5.08 | 2 | VV | 0.046 | 0.076 |
| 33 | 5.048 | 12.74 | 2 | VV | 0.082 | 0.191 |
| 34 | 5.199 | 6.89 | 2 | VV | 0.054 | 0.103 |
| 35 | 5.250 | 4.12 | 1 | VV | 0.043 | 0.062 |
| 36 | 5.351 | 731.57 | 493 | VV | 0.024 | 10.984 |

Area Percent Report

Data File Name : W:\GC_45\GC 45 DATA\2014\141030\14103064.D

Page Number : 2

Operator : 682 Vial Number : Vial 64

Instrument : GC 45 Injection Number : 1
Sample Name : 14-10-2131-3 Sequence Line : 64

Instrument Method: C:\CHEM32\1\METHODS\ ->

Acquired on : 31 Oct 14 09:59 am

Report Created on: 04 Nov 14 02:49 pm Analysis Method : 8015B.MTH

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Total area = 6660.03

Data File Name : W:\GC 45\GC 45 DATA\2014\141030\14103064.D

Page Number : 3

Operator : 682 Vial Number : Vial 64

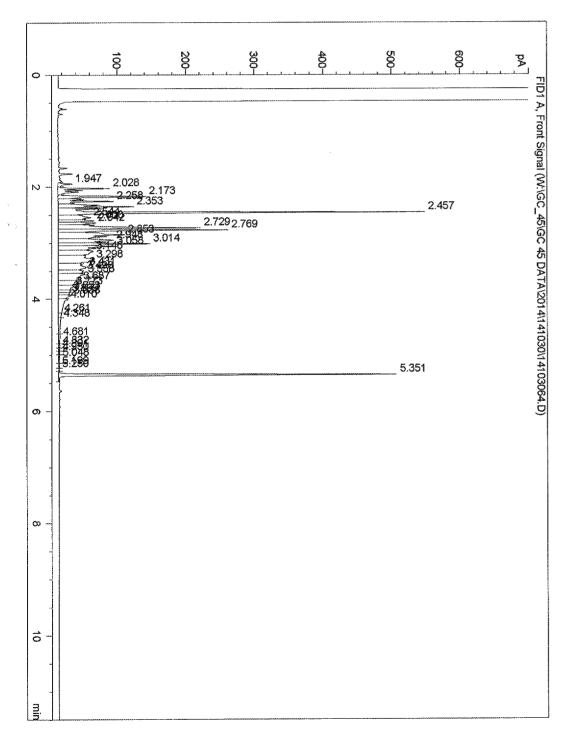
Instrument : GC 45 Injection Number : 1 Sample Name : 14-10-2131-3 Sequence Line : 64

Instrument Method: C:\CHEM32\1\METHODS\ ->

Acquired on : 31 Oct 14 09:59 am

Report Created on: 04 Nov 14 02:49 pm Analysis Method : 8015B.MTH

Software Revision: Rev. B.03.02 [341] Copyright © Agilent Technologies



RAW DATA SHEET FOR METHOD: EPA 8015B (M)

WORK ORDER: 14-10-2131

GC 45

EPA 3510C EXTRACTION:

2014-10-30 00:00 D/T EXTRACTED:

682 ANALYZED BY:

2014-10-31 10:41 D/T ANALYZED:

628 **REVIEWED BY:**

2014-11-04 17:13 D/T REVIEWED:

DATA FILE:

INSTRUMENT:

W:\GC 45\GC 45 DATA\2014\141030\14103065.D\14103065

4

UNITS:

CLIENT SAMPLE NUMBER: ES115

LCS/MB BATCH: MS/MSD BATCH: 141030B06

SAMPLE VOLUME / WEIGHT:

DEFAULT: 500.00 ml / ACTUAL: 500.00 ml

DEFAULT: 5.00 ml / ACTUAL: 2.50 ml FINAL VOLUME / WEIGHT:

ADJUSTMENT RATIO TO PF: 0.50

ug/L

Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are **COMMENT:**

qualified with a "J" flag. TPH as Diesel is quantified in the carbon range C10-C28.

QUAL LOQ CONC DL LOD INI. CONC COMPOUND DF

25 b 2030 11 12 406000 1.00 TPH as Diesel

Data File Name : W:\GC_45\GC 45 DATA\2014\141030\14103065.D

Vial Number : Vial 65

Page Number : 1
Operator : 682
Instrument : GC 45 Injection Number: 1 Sequence Line : 65 Sample Name : 14-10-2131-4

Instrument Method: C:\CHEM32\1\METHODS\

Acquired on : 31 Oct 14 10:41 am

Report Created on: 04 Nov 14 02:50 pm Analysis Method : 8015B.MTH

Software Revision: Rev. B.03.02 [341] Copyright © Agilent Technologies

Sig. 1 in W:\GC 45\GC 45 DATA\2014\141030\ ->

| Pk | in W:\GC_45\G Ret Time | C 45 DATA\2014 Area | Height | Peak | Width | Response % |
|----|---------------------------|------------------------|--------|------|-------|------------|
| 1 | 1.946 | 36.71 | | VV | 0.026 | 0.555 |
| 2 | 2.028 | 169.15 | | VV | 0.031 | 2.559 |
| 3 | 2.173 | 226.98 | 127 | VV | 0.028 | 3.434 |
| 4 | 2.222 | 61.36 | 36 | VV | 0.025 | 0.928 |
| 5 | 2.257 | 128.76 | 81 | VV | 0.024 | 1.948 |
| 6 | 2.352 | 231.89 | 109 | VV | 0.030 | 3.508 |
| 7 | 2.457 | 829.85 | 551 | VV | 0.023 | 12.555 |
| 8 | 2.542 | 189.89 | 47 | VV | 0.055 | 2.873 |
| 9 | 2.600 | 104.99 | 51 | VV | 0.030 | 1.588 |
| 10 | 2.641 | 115.18 | 56 | VV | 0.030 | 1.743 |
| 11 | 2.730 | 449.33 | 212 | VV | 0.030 | 6.798 |
| 12 | 2.770 | 375.36 | 260 | VV | 0.022 | 5.679 |
| 13 | 2.852 | 404.48 | 91 | VV | 0.057 | 6.119 |
| 14 | 2.947 | 210.06 | 78 | VV | 0.041 | 3.178 |
| 15 | 3.014 | 316.74 | 136 | VV | 0.032 | 4.792 |
| 16 | 3.059 | 267.38 | 81 | VV | 0.044 | 4.045 |
| 17 | 3.149 | 231.23 | | VV | 0.063 | 3.498 |
| 18 | 3.296 | 424.16 | | VV | 0.113 | 6.417 |
| 19 | 3.431 | 166.71 | | VV | 0.053 | 2.522 |
| 20 | 3.502 | 122.86 | | VV | 0.047 | 1.859 |
| 21 | 3.561 | 218.09 | | VV | 0.075 | 3.299 |
| 22 | 3.690 | 126.19 | | VV | 0.054 | 1.909 |
| 23 | 3.783 | 94.82 | | VV | 0.059 | 1.434 |
| 24 | 3.861 | 40.17 | | VV | 0.033 | 0.608 |
| 25 | 3.907 | 37.37 | | VV | 0.039 | 0.565 |
| 26 | 3.950 | 52.52 | | VV | 0.043 | 0.795 |
| 27 | 4.023 | 152.63 | | VV | 0.159 | 2.309 |
| 28 | 4.603 | 7.19 | | VV | 0.051 | 0.109 |
| 29 | 4.719 | 16.45 | | VV | 0.095 | 0.249 |
| 30 | 4.871 | 7.28 | | VV | 0.051 | 0.110 |
| 31 | 4.931 | 5.14 | | VV | 0.054 | 0.078 |
| 32 | 4.994 | 4.19 | | VV | 0.043 | 0.063 |
| 33 | 5.096 | 12.49 | | VV | 0.083 | 0.189 |
| 34 | 5.252 | 5.63 | | VV | 0.045 | 0.085 |
| 35 | 5.313 | 3.60 | | VV | 0.044 | 0.054 |
| 36 | 5.408 | 762.98 | 520 | VV | 0.023 | 11.543 |

Area Percent Report

Data File Name : W:\GC_45\GC 45 DATA\2014\141030\14103065.D

Vial Number : Vial 65

Page Number : 2
Operator : 682
Instrument : GC : GC 45 Injection Number: 1 Sample Name : 14-10-2131-4 Sequence Line : 65

Instrument Method: C:\CHEM32\1\METHODS\ ->

Acquired on : 31 Oct 14 10:41 am

Report Created on: 04 Nov 14 02:50 pm Analysis Method : 8015B.MTH

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Total area = 6609.83

Data File Name : W:\GC_45\GC 45 DATA\2014\141030\14103065.D

Page Number : 3

Operator : 682 Vial Number : Vial 65

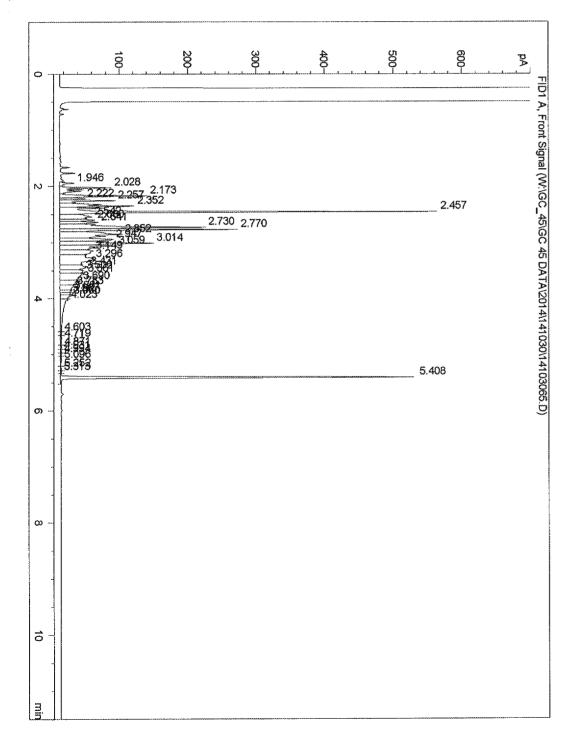
Instrument : GC 45 Injection Number : 1 Sample Name : 14-10-2131-4 Sequence Line : 65

Instrument Method: C:\CHEM32\1\METHODS\ ->

Acquired on : 31 Oct 14 10:41 am

Report Created on: 04 Nov 14 02:50 pm Analysis Method : 8015B.MTH

Software Revision: Rev. B.03.02 [341] Copyright © Agilent Technologies



RAW DATA SHEET FOR METHOD: EPA 8015B (M)

WORK ORDER:

D/T EXTRACTED:

INSTRUMENT:

EXTRACTION:

14-10-2131

GC 45

EPA 3510C

2014-10-30 00:00

ANALYZED BY:

682

2014-10-31 10:58 D/T ANALYZED:

REVIEWED BY: 628

D/T REVIEWED: 2014-11-04 17:13

DATA FILE:

W:\GC 45\GC 45 DATA\2014\141030\14103066.D\14103066

5

CLIENT SAMPLE NUMBER: ES116

LCS/MB BATCH:

141030B06 141030S06 SAMPLE VOLUME / WEIGHT:

DEFAULT: 500.00 ml / ACTUAL: 500.00 ml

DEFAULT: 5.00 ml / ACTUAL: 2.50 ml

MS/MSD BATCH: UNITS:

ug/L

FINAL VOLUME / WEIGHT: ADJUSTMENT RATIO TO PF:

11

COMMENT:

Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are

qualified with a "J" flag. TPH as Diesel is quantified in the carbon range C10-C28.

COMPOUND TPH as Diesel INI. CONC 16000

<u>DF</u> 1.00 CONC 79.9

LOD DL 12

LOQ

25

QUAL

b

Data File Name : W:\GC_45\GC 45 DATA\2014\141030\14103066.D

Page Number : 1

Vial Number : Vial 66

Operator : 682 Instrument : GC : GC 45 Injection Number: 1 Sequence Line : 66 Sample Name : 14-10-2131-5

> Instrument Method: C:\CHEM32\1\METHODS\ -->

Acquired on : 31 Oct 14 10:58 am

Report Created on: 04 Nov 14 02:51 pm Analysis Method : 8015B.MTH

Software Revision: Rev. B.03.02 [341] Copyright © Agilent Technologies

Sig. 1 in W:\GC 45\GC 45 DATA\2014\141030\ \rightarrow

| Pk | Ret Time | Area | - | | | - |
|----|----------|--------|-----|----|-------|--------|
| 1 | | | | | | . |
| 1 | 2.219 | 0.45 | 0 | VV | 0.038 | 0.060 |
| 2 | 2.270 | 0.27 | 0 | VV | 0.031 | 0.036 |
| 3 | 2.457 | 0.84 | 0 | VV | 0.065 | 0.113 |
| 4 | 2.634 | 1.87 | 1 | VV | 0.049 | 0.252 |
| 5 | 2.750 | 4.97 | 1 | VV | 0.064 | 0.669 |
| 6 | 2.837 | 1.49 | 1 | VV | 0.037 | 0.200 |
| 7 | 2.884 | 1.51 | 1 | VV | 0.032 | 0.204 |
| 8 | 2.950 | 2.03 | 1 | VV | 0.045 | 0.273 |
| 9 | 3.169 | 12.58 | 2 | VV | 0.078 | 1.695 |
| 10 | 3.306 | 10.78 | 2 | VV | 0.070 | 1.452 |
| 11 | 3.372 | 8.28 | 2 | VV | 0.058 | 1.115 |
| 12 | 3.436 | 8.46 | 3 | VV | 0.046 | 1.139 |
| 13 | 3.511 | 26.81 | 3 | VV | 0.117 | 3.611 |
| 14 | 3.701 | 20.15 | 2 | VV | 0.122 | 2.714 |
| 15 | 3.846 | 7.00 | 2 | VV | 0.051 | 0.943 |
| 16 | 3.937 | 5.92 | 2 | VV | 0.052 | 0.797 |
| 17 | 4.012 | 7.26 | 3 | VV | 0.036 | 0.978 |
| 18 | 4.111 | 8.62 | 2 | VV | 0.075 | 1.161 |
| 19 | 4.228 | 13.65 | 2 | VV | 0.096 | 1.839 |
| 20 | 4.387 | 20.35 | 2 | VV | 0.137 | 2.740 |
| 21 | 4.560 | 13.57 | 2 | VV | 0.100 | 1.828 |
| 22 | 4.662 | 12.72 | 2 | VV | 0.090 | 1.714 |
| 23 | 4.826 | 7.22 | 2 | VV | 0.064 | 0.972 |
| 24 | 4,889 | 3.85 | 1 | VV | 0.042 | 0.518 |
| 25 | 4.946 | 6.18 | 1 | VV | 0.061 | 0.832 |
| 26 | 5.043 | 8.15 | 1 | VV | 0.074 | 1.097 |
| 27 | 5.193 | 3.65 | 1 | VV | 0.041 | 0.492 |
| 28 | 5.249 | 2.50 | 1 | VV | 0.048 | 0.337 |
| 29 | 5.344 | 521.35 | 374 | VV | 0.022 | 70.220 |

Total area = 742.45

Data File Name : W:\GC_45\GC 45 DATA\2014\141030\14103066.D

Page Number : 2

: 682 Operator

: GC 45 Instrument

Sample Name

: 14-10-2131-5

Injection Number: 1

: 66 Sequence Line

Vial Number : Vial 66

Instrument Method: C:\CHEM32\1\METHODS\

Acquired on

: 31 Oct 14

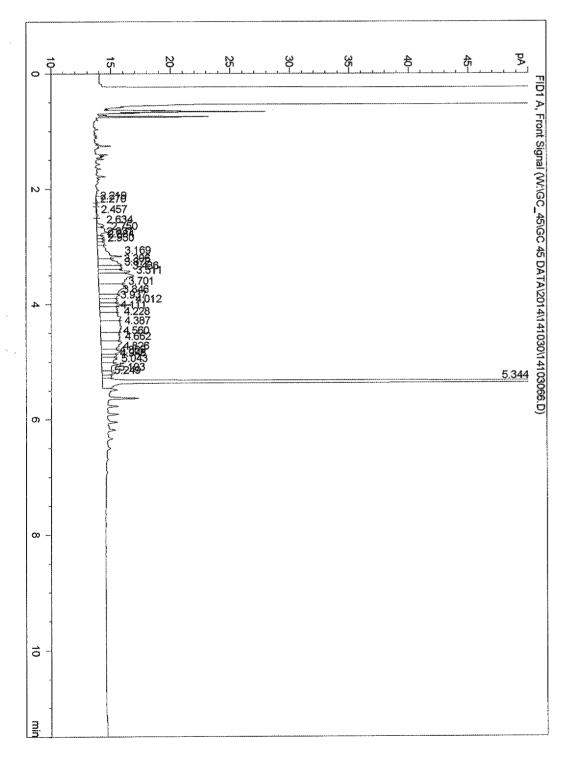
Report Created on: 04 Nov 14

10:58 am

02:51 pm

Analysis Method: 8015B.MTH

Software Revision: Rev. B.03.02 [341] Copyright © Agilent Technologies



Data File Name : W:\GC 45\GC 45 DATA\2014\141030\14103051.D

Page Number : 3

Operator : 682 Vial Number : Vial 51

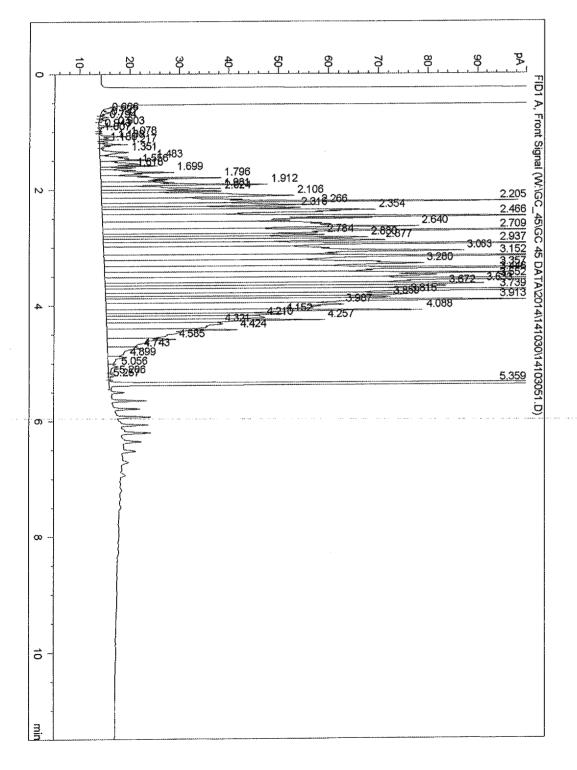
Instrument : GC 45 Injection Number : 1 Sample Name : D400 C28 50 L080814D Sequence Line : 51

Instrument Method: C:\CHEM32\1\METHODS\ ->

Acquired on : 31 Oct 14 06:11 am

Report Created on: 04 Nov 14 01:18 pm Analysis Method : 8015B.MTH

Software Revision: Rev. B.03.02 [341] Copyright © Agilent Technologies





Calscience



WORK ORDER NUMBER: 14-10-2234

The difference is service



AIR SOIL WATER | MARINE CHEMISTRY

Analytical Report For

Client: Environmental Science International, Inc.

Client Project Name: Red Hill LTM 112066

Attention: Jeff Hattemer

354 Uluniu Street, Suite 304 Kailua, HI 96734-2500

Rand Odl

Approved for release on 11/05/2014 by: Richard Villafania Project Manager



Emailyour PM)

Resulting

Eurofins Calscience, Inc. (Calscience) certifies that the test results provided in this report meet all NELAC requirements for parameters for which accreditation is required or available. Any exceptions to NELAC requirements are noted in the case narrative. The original report of subcontracted analyses, if any, is attached to this report. The results in this report are limited to the sample(s) tested and any reproduction thereof must be made in its entirety. The client or recipient of this report is specifically prohibited from making material changes to said report and, to the extent that such changes are made, Calscience is not responsible, legally or otherwise. The client or recipient agrees to indemnify Calscience for any defense to any litigation which may arise.

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NELAP DI 002/00A I ACLASS DODIELAP ID ADE 1984 (ISDREC 17025/2025) I OSDLAG DI 10109 I SCAQMO DI POLADERS



Contents

| Client Project Name: | Red Hill LTM 112066 |
|----------------------|---------------------|
| Work Order Number: | 14-10-2234 |

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Work Order Narrative

Work Order: 14-10-2234 Page 1 of 1

Condition Upon Receipt:

Samples were received under Chain-of-Custody (COC) on 10/29/14. They were assigned to Work Order 14-10-2234.

Unless otherwise noted on the Sample Receiving forms all samples were received in good condition and within the recommended EPA temperature criteria for the methods noted on the COC. The COC and Sample Receiving Documents are integral elements of the analytical report and are presented at the back of the report.

Holding Times:

All samples were analyzed within prescribed holding times (HT) and/or in accordance with the Calscience Sample Acceptance Policy unless otherwise noted in the analytical report and/or comprehensive case narrative, if required.

Any parameter identified in 40CFR Part 136.3 Table II that is designated as "analyze immediately" with a holding time of <= 15 minutes (40CFR-136.3 Table II, footnote 4), is considered a "field" test and the reported results will be qualified as being received outside of the stated holding time unless received at the laboratory within 15 minutes of the collection time.

Quality Control:

All quality control parameters (QC) were within established control limits except where noted in the QC summary forms or described further within this report.

Additional Comments:

Air - Sorbent-extracted air methods (EPA TO-4A, EPA TO-10, EPA TO-13A, EPA TO-17): Analytical results are converted from mass/sample basis to mass/volume basis using client-supplied air volumes.

New York NELAP air certification does not certify for all reported methods and analytes, reference the accredited items here: http://www.calscience.com/PDF/New_York.pdf

Solid - Unless otherwise indicated, solid sample data is reported on a wet weight basis, not corrected for % moisture. All QC results are always reported on a wet weight basis.

Subcontractor Information:

Unless otherwise noted below (or on the subcontract form), no samples were subcontracted.



10/29/14 Date Received: Environmental Science International, Inc. Work Order: 14-10-2234 354 Uluniu Street, Suite 304 Kailua, HI 96734-2500 Preparation: **EPA 3510C** Method: EPA 8015B (M) Units: ug/L

Project: Red Hill LTM 112066 Page 1 of 1 Lab Sample Client Sample Number Date/Time Matrix Instrument Date Date/Time QC Batch ID Number Collected Prepared Analyzed

ES117 14-10-2234-2-G 10/28/14 Aqueous GC 45 10/30/14 10/31/14 141030B06

Comment(s): - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

- TPH as Diesel is quantified in the carbon range C10-C28.

Parameter Result DL LOD LOQ DF Qualifiers TPH as Diesel 22 11 12 25 1.00 HD,J

Surrogate Rec. (%) **Control Limits** Qualifiers

51-141 n-Octacosane 85

ES118 10/28/14 14:30 14-10-2234-3-G 10/30/14 141030B06 10/31/14 Aqueous GC 45 11:35 - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

Comment(s):

- TPH as Diesel is quantified in the carbon range C10-C28.

<u>LOD</u> <u>LOQ</u> Qualifiers <u>Parameter</u> Result <u>DF</u> TPH as Diesel 16 11 12 25 1.00 HD,J

Control Limits Qualifiers Surrogate Rec. (%)

n-Octacosane 70 51-141

| Method Blank | 099- | 15-516-206 N/ | A Aqı | ueous GC 45 | 10/30/14 | 10/31/14 07:03 | 141030B06 |
|----------------------------------|-----------------------------------|-------------------|--------------------------|-------------------|------------------|--------------------|-------------------|
| Comment(s): | - Results were evaluated to the M | DL (DL), concentr | ations >= to the I | MDL (DL) but < RL | (LOQ), if found, | are qualified with | a "J" flag. |
| <u>Parameter</u> | | <u>Result</u> | <u>DL</u> | LOD | LOQ | <u>DF</u> | <u>Qualifiers</u> |
| TPH as Diesel | | <12 | 11 | 12 | 25 | 1.00 | U |
| <u>Surrogate</u> n-Octacosane | | Rec. (%) | Control Limits 51-141 | <u>Qualifiers</u> | | | |



Environmental Science International, Inc.

Work Order:

10/29/14

354 Uluniu Street, Suite 304 Kailua, HI 96734-2500

Preparation:

Date Received:

EPA 3005A Filt.

Method:

EPA 6020

Units:

ug/L

14-10-2234

Project: Red Hill LTM 112066

Page 1 of 1

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|---|-----------------------|------------------------|------------|------------------|-------------------|-----------------------|-------------|
| ES118 | 14-10-2234-3-J | 10/28/14 14:30 | Aqueous | ICP/MS 04 | 10/31/14 | 11/03/14 15:10 | 141031L04F |
| Comment(s): - Results were evaluated to | to the MDL (DL), cond | centrations >= t | the MDL (D | DL) but < RL (LO | Q), if found, are | qualified with a | "J" flag. |

 Parameter
 Result
 DL
 LOD
 LOQ
 DF
 Qualifiers

 Lead
 <0.200</td>
 0.0898
 0.200
 1.00
 1.00
 U

| Method Blank | 099-14-497-10 | 4 N/A | Aqueous ICP/ | MS 04 10/31/ | 14 11/03/14 14:48 | 1 141031L04F |
|------------------|---|----------------------|--------------------|---------------------|----------------------|-------------------|
| Comment(s): | - Results were evaluated to the MDL (DL), o | concentrations >= to | o the MDL (DL) but | < RL (LOQ), if four | nd, are qualified w | vith a "J" flag. |
| <u>Parameter</u> | Result | <u>DL</u> | LOD | LOQ | <u>DF</u> | <u>Qualifiers</u> |
| Lead | <0.200 | 0.0898 | 0.200 | 1.00 | 1.00 | U |





Environmental Science International, Inc.

Kailua, HI 96734-2500

Date Received:

10/29/14 14-10-2234

354 Uluniu Street, Suite 304

Work Order: Preparation:

EPA 3510C

Method:

EPA 8270C SIM PAHs

Units:

ug/L

Project: Red Hill LTM 112066

Page 1 of 3

| Client Sample Number | Lab Sample Number | Date/Time I Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|--------------------------------------|------------------------|--------------------------|-------------------|-----------------|------------------|-----------------------|-------------------|
| ES117 | 14-10-2234-2-1 | 10/28/14 12:00 | Aqueous | GC/MS AAA | 10/30/14 | 11/03/14 17:03 | 141030L08 |
| Comment(s): - Results were evaluated | d to the MDL (DL), con | centrations >= to th | ne MDL (DL |) but < RL (LOC | Q), if found, a | are qualified with | a "J" flag. |
| <u>Parameter</u> | Result | <u>DL</u> | LOD | LOC | 3 | <u>DF</u> | <u>Qualifiers</u> |
| Naphthalene | <0.049 | 0.033 | 0.049 | 0.19 |) | 1.00 | U |
| 2-Methylnaphthalene | < 0.049 | 0.045 | 0.049 | 0.19 |) | 1.00 | U |
| 1-Methylnaphthalene | <0.097 | 0.050 | 0.097 | 0.19 |) | 1.00 | U |
| Acenaphthylene | < 0.049 | 0.043 | 0.049 | 0.19 |) | 1.00 | U |
| Acenaphthene | <0.049 | 0.026 | 0.049 | 0.19 |) | 1.00 | U |
| Fluorene | < 0.049 | 0.041 | 0.049 | 0.19 |) | 1.00 | U |
| Phenanthrene | <0.049 | 0.026 | 0.049 | 0.19 |) | 1.00 | U |
| Anthracene | < 0.049 | 0.028 | 0.049 | 0.19 |) | 1.00 | U |
| Fluoranthene | < 0.049 | 0.045 | 0.049 | 0.19 |) | 1.00 | U |
| Pyrene | < 0.049 | 0.020 | 0.049 | 0.19 |) | 1.00 | U |
| Benzo (a) Anthracene | < 0.049 | 0.032 | 0.049 | 0.19 |) | 1.00 | U |
| Chrysene | < 0.049 | 0.024 | 0.049 | 0.19 |) | 1.00 | U |
| Benzo (k) Fluoranthene | < 0.049 | 0.030 | 0.049 | 0.19 |) | 1.00 | U |
| Benzo (b) Fluoranthene | < 0.049 | 0.017 | 0.049 | 0.19 |) | 1.00 | U |
| Benzo (a) Pyrene | < 0.049 | 0.022 | 0.049 | 0.19 |) | 1.00 | U |
| Indeno (1,2,3-c,d) Pyrene | < 0.049 | 0.021 | 0.049 | 0.19 |) | 1.00 | U |
| Dibenz (a,h) Anthracene | < 0.049 | 0.046 | 0.049 | 0.19 |) | 1.00 | U |
| Benzo (g,h,i) Perylene | <0.097 | 0.080 | 0.097 | 0.19 |) | 1.00 | U |
| Surrogate | Rec. (%) | Control Limit | ts <u>Qualifi</u> | <u>ers</u> | | | |
| Nitrobenzene-d5 | 61 | 28-139 | | | | | |
| 2-Fluorobiphenyl | 81 | 33-144 | | | | | |
| p-Terphenyl-d14 | 83 | 23-160 | | | | | |



Environmental Science International, Inc.

354 Uluniu Street, Suite 304 Kailua, HI 96734-2500 Date Received:

10/29/14 14-10-2234

Work Order: Preparation:

EPA 3510C

Method:

EPA 8270C SIM PAHs

Units:

ug/L

Project: Red Hill LTM 112066

Page 2 of 3

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | | Date/Time QC Batch ID Analyzed |
|--------------------------------------|----------------------|------------------------|-------------|-----------------|----------------------|-----------------------------------|
| ES118 | 14-10-2234-3-1 | 10/28/14 14:30 | Aqueous | GC/MS AAA | | 11/03/14 141030L08 17:26 |
| Comment(s): - Results were evaluated | to the MDL (DL), con | centrations >= to th | he MDL (DL) |) but < RL (LOC | i), if found, are qu | alified with a "J" flag. |
| <u>Parameter</u> | Result | <u>DL</u> | LOD | LOC | <u>DF</u> | <u>Qualifiers</u> |
| Naphthalene | <0.048 | 0.033 | 0.048 | 0.19 | 1.00 |) U |
| 2-Methylnaphthalene | <0.048 | 0.045 | 0.048 | 0.19 | 1.00 |) U |
| 1-Methylnaphthalene | <0.096 | 0.050 | 0.096 | 0.19 | 1.00 |) U |
| Acenaphthylene | <0.048 | 0.043 | 0.048 | 0.19 | 1.00 |) U |
| Acenaphthene | <0.048 | 0.026 | 0.048 | 0.19 | 1.00 |) U |
| Fluorene | <0.048 | 0.041 | 0.048 | 0.19 | 1.00 |) U |
| Phenanthrene | <0.048 | 0.026 | 0.048 | 0.19 | 1.00 |) U |
| Anthracene | <0.048 | 0.028 | 0.048 | 0.19 | 1.00 |) U |
| Fluoranthene | <0.048 | 0.045 | 0.048 | 0.19 | 1.00 |) U |
| Pyrene | <0.048 | 0.020 | 0.048 | 0.19 | 1.00 |) U |
| Benzo (a) Anthracene | <0.048 | 0.031 | 0.048 | 0.19 | 1.00 |) U |
| Chrysene | <0.048 | 0.024 | 0.048 | 0.19 | 1.00 |) U |
| Benzo (k) Fluoranthene | <0.048 | 0.030 | 0.048 | 0.19 | 1.00 |) U |
| Benzo (b) Fluoranthene | <0.048 | 0.017 | 0.048 | 0.19 | 1.00 |) U |
| Benzo (a) Pyrene | <0.048 | 0.021 | 0.048 | 0.19 | 1.00 |) U |
| Indeno (1,2,3-c,d) Pyrene | <0.048 | 0.020 | 0.048 | 0.19 | 1.00 |) U |
| Dibenz (a,h) Anthracene | <0.048 | 0.046 | 0.048 | 0.19 | 1.00 |) U |
| Benzo (g,h,i) Perylene | <0.096 | 0.079 | 0.096 | 0.19 | 1.00 |) U |
| Surrogate | Rec. (%) | Control Limi | its Qualifi | ers | | |
| Nitrobenzene-d5 | 62 | 28-139 | | | | |
| 2-Fluorobiphenyl | 88 | 33-144 | | | | |
| p-Terphenyl-d14 | 94 | 23-160 | | | | |
| • • • | | | | | | |



Environmental Science International, Inc.

Kailua, HI 96734-2500

Date Received:

10/29/14 14-10-2234

354 Uluniu Street, Suite 304

Work Order: Preparation:

EPA 3510C

Method:

EPA 8270C SIM PAHs

Units:

ug/L

Project: Red Hill LTM 112066

Page 3 of 3

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|--------------------------------------|----------------------|------------------------|-------------|-----------------|--------------------|-----------------------|-------------|
| Method Blank | 099-15-148-67 | N/A | Aqueous | GC/MS AAA | 10/30/14 | 10/31/14 17:17 | 141030L08 |
| Comment(s): - Results were evaluated | to the MDL (DL), con | ncentrations >= to the | he MDL (DL) |) but < RL (LOC |), if found, are o | ualified with a | "J" flag. |
| <u>Parameter</u> | <u>Result</u> | <u>DL</u> | LOD | LOC | <u>DF</u> | : | Qualifiers |
| Naphthalene | <0.050 | 0.034 | 0.050 | 0.20 | 1.0 | 00 | U |
| 2-Methylnaphthalene | <0.050 | 0.046 | 0.050 | 0.20 | 1.0 | 00 | U |
| 1-Methylnaphthalene | <0.10 | 0.052 | 0.10 | 0.20 | 1.0 | 00 | U |
| Acenaphthylene | <0.050 | 0.045 | 0.050 | 0.20 | 1.0 | 00 | U |
| Acenaphthene | <0.050 | 0.027 | 0.050 | 0.20 | 1.0 | 00 | U |
| Fluorene | <0.050 | 0.043 | 0.050 | 0.20 | 1.0 | 00 | U |
| Phenanthrene | <0.050 | 0.027 | 0.050 | 0.20 | 1.0 | 00 | U |
| Anthracene | <0.050 | 0.029 | 0.050 | 0.20 | 1.0 | 00 | U |
| Fluoranthene | <0.050 | 0.047 | 0.050 | 0.20 | 1.0 | 00 | U |
| Pyrene | <0.050 | 0.020 | 0.050 | 0.20 | 1.0 | 00 | U |
| Benzo (a) Anthracene | <0.050 | 0.033 | 0.050 | 0.20 | 1.0 | 00 | U |
| Chrysene | <0.050 | 0.025 | 0.050 | 0.20 | 1.0 | 00 | U |
| Benzo (k) Fluoranthene | <0.050 | 0.031 | 0.050 | 0.20 | 1.0 | 00 | U |
| Benzo (b) Fluoranthene | <0.050 | 0.018 | 0.050 | 0.20 | 1.0 | 00 | U |
| Benzo (a) Pyrene | <0.050 | 0.022 | 0.050 | 0.20 | 1.0 | 00 | U |
| Indeno (1,2,3-c,d) Pyrene | <0.050 | 0.021 | 0.050 | 0.20 | 1.0 | 00 | U |
| Dibenz (a,h) Anthracene | <0.050 | 0.048 | 0.050 | 0.20 | 1.0 | 00 | U |
| Benzo (g,h,i) Perylene | <0.10 | 0.082 | 0.10 | 0.20 | 1.0 | 00 | U |
| Surrogate | Rec. (%) | Control Limi | its Qualifi | ers | | | |
| Nitrobenzene-d5 | 53 | 28-139 | | | | | |
| 2-Fluorobiphenyl | 68 | 33-144 | | | | | |
| p-Terphenyl-d14 | 72 | 23-160 | | | | | |
| L L | · | | | | | | |



Environmental Science International, Inc.

354 Uluniu Street, Suite 304 Kailua, HI 96734-2500 Date Received:

10/29/14 14-10-2234

Work Order: Preparation:

EPA 5030C

Method:

GC/MS / EPA 8260B

Units:

ug/L

Project: Red Hill LTM 112066

Page 1 of 8

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Date/Time Prepared Analyzed | QC Batch ID |
|---|----------------------|------------------------|----------------|------------------|-------------------------------------|-------------------|
| ES TRIP | 14-10-2234-1-A | 10/28/14 11:00 | Aqueous | GC/MS OO | 10/31/14 11/01/14 01:14 | 141031L047 |
| Comment(s): - Results were evaluated to | o the MDL (DL), con | centrations >= | to the MDL (DI | _) but < RL (LOC |), if found, are qualified with | ı a "J" flag. |
| <u>Parameter</u> | <u>Result</u> | <u>DL</u> | <u>LOD</u> | LOC | <u>DF</u> | <u>Qualifiers</u> |
| Acetone | <10 | 6.0 | 10 | 20 | 1.00 | U,ICH |
| Benzene | <0.50 | 0.14 | 0.50 | 1.0 | 1.00 | U |
| Bromodichloromethane | <0.50 | 0.21 | 0.50 | 5.0 | 1.00 | U |
| Bromoform | <1.0 | 0.50 | 1.0 | 10 | 1.00 | U |
| Bromomethane | <5.0 | 3.9 | 5.0 | 20 | 1.00 | U,IJ |
| 2-Butanone | <5.0 | 2.2 | 5.0 | 10 | 1.00 | U |
| Carbon Tetrachloride | <0.50 | 0.23 | 0.50 | 1.0 | 1.00 | U |
| Chlorobenzene | <0.50 | 0.17 | 0.50 | 5.0 | 1.00 | U |
| Chloroethane | <5.0 | 2.3 | 5.0 | 10 | 1.00 | U |
| Chloroform | <0.50 | 0.46 | 0.50 | 5.0 | 1.00 | U |
| Chloromethane | <2.0 | 1.8 | 2.0 | 10 | 1.00 | U |
| Dibromochloromethane | <0.50 | 0.25 | 0.50 | 1.0 | 1.00 | U |
| 1,2-Dibromo-3-Chloropropane | <2.0 | 1.2 | 2.0 | 10 | 1.00 | U |
| 1,2-Dibromoethane | <0.50 | 0.36 | 0.50 | 1.0 | 1.00 | U |
| 1,2-Dichlorobenzene | <0.50 | 0.46 | 0.50 | 1.0 | 1.00 | U |
| 1,3-Dichlorobenzene | <0.50 | 0.40 | 0.50 | 1.0 | 1.00 | U |
| 1,4-Dichlorobenzene | 0.43 | 0.43 | 0.50 | 1.0 | 1.00 | J |
| 1,1-Dichloroethane | <0.50 | 0.28 | 0.50 | 5.0 | 1.00 | U |
| 1,2-Dichloroethane | <0.50 | 0.24 | 0.50 | 1.0 | 1.00 | U |
| 1,1-Dichloroethene | <0.50 | 0.43 | 0.50 | 1.0 | 1.00 | U |
| c-1,2-Dichloroethene | <0.50 | 0.48 | 0.50 | 1.0 | 1.00 | U |
| t-1,2-Dichloroethene | <0.50 | 0.37 | 0.50 | 1.0 | 1.00 | U |
| 1,2-Dichloropropane | <0.50 | 0.42 | 0.50 | 5.0 | 1.00 | U |
| c-1,3-Dichloropropene | <0.50 | 0.25 | 0.50 | 1.0 | 1.00 | U |
| t-1,3-Dichloropropene | <0.50 | 0.25 | 0.50 | 1.0 | 1.00 | U |
| Ethylbenzene | <0.50 | 0.14 | 0.50 | 1.0 | 1.00 | U |
| Methylene Chloride | <1.0 | 0.64 | 1.0 | 5.0 | 1.00 | U |
| 4-Methyl-2-Pentanone | <5.0 | 4.4 | 5.0 | 10 | 1.00 | U |
| Styrene | <0.50 | 0.17 | 0.50 | 1.0 | 1.00 | U |
| 1,1,1,2-Tetrachloroethane | <0.50 | 0.40 | 0.50 | 1.0 | 1.00 | U |
| 1,1,2,2-Tetrachloroethane | <0.50 | 0.41 | 0.50 | 1.0 | 1.00 | U |
| Tetrachloroethene | <0.50 | 0.39 | 0.50 | 5.0 | 1.00 | U |
| Toluene | <0.50 | 0.24 | 0.50 | 1.0 | 1.00 | U |
| 1,2,4-Trichlorobenzene | <1.0 | 0.50 | 1.0 | 5.0 | 1.00 | U |
| 1.1.1-Trichloroethane | <0.50 | 0.30 | 0.50 | 5.0 | 1.00 | U |
| Hexachloro-1,3-Butadiene | <0.50 | 0.32 | 0.50 | 1.0 | 1.00 | U |



Environmental Science International, Inc. 354 Uluniu Street, Suite 304

Kailua, HI 96734-2500

Date Received: Work Order:

Preparation: Method: Units: 10/29/14

14-10-2234 EPA 5030C

GC/MS / EPA 8260B

ug/L

Project: Red Hill LTM 112066

Page 2 of 8

| Parameter | Result | <u>DL</u> | LOD | LOQ | <u>DF</u> | Qualifiers |
|-----------------------------|----------|----------------|------------|-----|-----------|------------|
| 1,1,2-Trichloroethane | <0.50 | 0.38 | 0.50 | 1.0 | 1.00 | U |
| Trichloroethene | <0.50 | 0.37 | 0.50 | 1.0 | 1.00 | U |
| 1,2,3-Trichloropropane | <1.0 | 0.64 | 1.0 | 5.0 | 1.00 | U |
| Vinyl Chloride | <0.50 | 0.30 | 0.50 | 1.0 | 1.00 | U |
| p/m-Xylene | <1.0 | 0.30 | 1.0 | 10 | 1.00 | U |
| o-Xylene | <0.50 | 0.23 | 0.50 | 1.0 | 1.00 | U |
| Methyl-t-Butyl Ether (MTBE) | <0.50 | 0.31 | 0.50 | 1.0 | 1.00 | U |
| Gasoline Range Organics | <30 | 26 | 30 | 50 | 1.00 | U |
| Surrogate | Rec. (%) | Control Limits | Qualifiers | | | |
| Dibromofluoromethane | 96 | 80-126 | | | | |
| 1,2-Dichloroethane-d4 | 100 | 80-134 | | | | |
| Toluene-d8 | 99 | 80-120 | | | | |
| Toluene-d8-TPPH | 95 | 88-112 | | | | |
| 1.4-Bromofluorobenzene | 97 | 80-120 | | | | |

10/29/14

14-10-2234

EPA 5030C



Analytical Report

Date Received:

Environmental Science International, Inc.

354 Uluniu Street, Suite 304 Kailua, HI 96734-2500

Work Order: Preparation: Method: GC/MS / EPA 8260B

Units: ug/L

Project: Red Hill LTM 112066 Page 3 of 8

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument Date Prepa | | Date/Time Analyzed | QC Batch ID |
|--|----------------------|------------------------|----------------|--------------------------|------------------|-----------------------|-------------------|
| ES117 | 14-10-2234-2-A | 10/28/14 12:00 | Aqueous | GC/MS OO | 10/31/14 | 11/01/14 02:36 | 141031L047 |
| Comment(s): - Results were evaluated t | o the MDL (DL), con | centrations >= t | to the MDL (DI | _) but < RL (LOC | (a), if found, a | re qualified with a | a "J" flag. |
| <u>Parameter</u> | <u>Result</u> | <u>DL</u> | <u>LOD</u> | LOC | Š | <u>DF</u> | <u>Qualifiers</u> |
| Acetone | <10 | 6.0 | 10 | 20 | | 1.00 | U,ICH |
| Benzene | <0.50 | 0.14 | 0.50 | 1.0 | | 1.00 | U |
| Bromodichloromethane | <0.50 | 0.21 | 0.50 | 5.0 | | 1.00 | U |
| Bromoform | <1.0 | 0.50 | 1.0 | 10 | | 1.00 | U |
| Bromomethane | <5.0 | 3.9 | 5.0 | 20 | | 1.00 | U,IJ |
| 2-Butanone | <5.0 | 2.2 | 5.0 | 10 | | 1.00 | U |
| Carbon Tetrachloride | <0.50 | 0.23 | 0.50 | 1.0 | | 1.00 | U |
| Chlorobenzene | <0.50 | 0.17 | 0.50 | 5.0 | | 1.00 | U |
| Chloroethane | <5.0 | 2.3 | 5.0 | 10 | | 1.00 | U |
| Chloroform | <0.50 | 0.46 | 0.50 | 5.0 | | 1.00 | U |
| Chloromethane | <2.0 | 1.8 | 2.0 | 10 | | 1.00 | U |
| Dibromochloromethane | <0.50 | 0.25 | 0.50 | 1.0 | | 1.00 | U |
| 1,2-Dibromo-3-Chloropropane | <2.0 | 1.2 | 2.0 | 10 | | 1.00 | U |
| 1,2-Dibromoethane | <0.50 | 0.36 | 0.50 | 1.0 | | 1.00 | U |
| 1,2-Dichlorobenzene | <0.50 | 0.46 | 0.50 | 1.0 | | 1.00 | U |
| 1,3-Dichlorobenzene | <0.50 | 0.40 | 0.50 | 1.0 | | 1.00 | U |
| 1,4-Dichlorobenzene | <0.50 | 0.43 | 0.50 | 1.0 | | 1.00 | U |
| 1,1-Dichloroethane | <0.50 | 0.28 | 0.50 | 5.0 | | 1.00 | U |
| 1,2-Dichloroethane | <0.50 | 0.24 | 0.50 | 1.0 | | 1.00 | U |
| 1,1-Dichloroethene | <0.50 | 0.43 | 0.50 | 1.0 | | 1.00 | U |
| c-1,2-Dichloroethene | <0.50 | 0.48 | 0.50 | 1.0 | | 1.00 | U |
| t-1,2-Dichloroethene | <0.50 | 0.37 | 0.50 | 1.0 | | 1.00 | U |
| 1,2-Dichloropropane | <0.50 | 0.42 | 0.50 | 5.0 | | 1.00 | U |
| c-1,3-Dichloropropene | <0.50 | 0.25 | 0.50 | 1.0 | | 1.00 | U |
| t-1,3-Dichloropropene | <0.50 | 0.25 | 0.50 | 1.0 | | 1.00 | U |
| Ethylbenzene | <0.50 | 0.14 | 0.50 | 1.0 | | 1.00 | U |
| Methylene Chloride | <1.0 | 0.64 | 1.0 | 5.0 | | 1.00 | U |
| 4-Methyl-2-Pentanone | <5.0 | 4.4 | 5.0 | 10 | | 1.00 | U |
| Styrene | <0.50 | 0.17 | 0.50 | 1.0 | | 1.00 | U |
| 1,1,2-Tetrachloroethane | <0.50 | 0.40 | 0.50 | 1.0 | | 1.00 | U |
| 1,1,2,2-Tetrachloroethane | <0.50 | 0.41 | 0.50 | 1.0 | | 1.00 | U |
| Tetrachloroethene | <0.50 | 0.39 | 0.50 | 5.0 | | 1.00 | U |
| Toluene | <0.50 | 0.24 | 0.50 | 1.0 | | 1.00 | U |
| 1,2,4-Trichlorobenzene | <1.0 | 0.50 | 1.0 | 5.0 | | 1.00 | U |
| 1,1,1-Trichloroethane | <0.50 | 0.30 | 0.50 | 5.0 | | 1.00 | U |
| Hexachloro-1,3-Butadiene | <0.50 | 0.32 | 0.50 | 1.0 | | 1.00 | U |



Environmental Science International, Inc. 354 Uluniu Street, Suite 304

Kailua, HI 96734-2500

Date Received:

Work Order: Preparation:

Method: Units: 10/29/14 14-10-2234

EPA 5030C

GC/MS / EPA 8260B

ug/L

Project: Red Hill LTM 112066

Page 4 of 8

| <u>Parameter</u> | Result | DL | LOD | LOQ | <u>DF</u> | Qualifiers |
|-----------------------------|----------|----------------|------------|-----|-----------|------------|
| 1,1,2-Trichloroethane | <0.50 | 0.38 | 0.50 | 1.0 | 1.00 | U |
| Trichloroethene | <0.50 | 0.37 | 0.50 | 1.0 | 1.00 | U |
| 1,2,3-Trichloropropane | <1.0 | 0.64 | 1.0 | 5.0 | 1.00 | U |
| Vinyl Chloride | <0.50 | 0.30 | 0.50 | 1.0 | 1.00 | U |
| p/m-Xylene | <1.0 | 0.30 | 1.0 | 10 | 1.00 | U |
| o-Xylene | <0.50 | 0.23 | 0.50 | 1.0 | 1.00 | U |
| Methyl-t-Butyl Ether (MTBE) | <0.50 | 0.31 | 0.50 | 1.0 | 1.00 | U |
| Gasoline Range Organics | <30 | 26 | 30 | 50 | 1.00 | U |
| Surrogate | Rec. (%) | Control Limits | Qualifiers | | | |
| Dibromofluoromethane | 96 | 80-126 | | | | |
| 1,2-Dichloroethane-d4 | 99 | 80-134 | | | | |
| Toluene-d8 | 99 | 80-120 | | | | |
| Toluene-d8-TPPH | 95 | 88-112 | | | | |
| 1,4-Bromofluorobenzene | 95 | 80-120 | | | | |



Environmental Science International, Inc.

Date Received: 10/29/14

354 Uluniu Street, Suite 304

Work Order: 14-10-2234

 Kailua, HI 96734-2500
 Preparation:
 EPA 5030C

 Method:
 GC/MS / EPA 8260B

Units: ug/L

Project: Red Hill LTM 112066 Page 5 of 8

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Date/Time Prepared Analyzed | QC Batch ID |
|--------------------------------------|------------------------|------------------------|---------------|------------------|-------------------------------------|---------------|
| ES118 | 14-10-2234-3-A | 10/28/14 14:30 | Aqueous | GC/MS OO | 10/31/14 11/01/14 03:04 | 141031L047 |
| Comment(s): - Results were evaluated | I to the MDL (DL), con | centrations >= t | o the MDL (DI | _) but < RL (LOG |), if found, are qualified wit | h a "J" flag. |
| <u>Parameter</u> | Result | <u>DL</u> | LOD | LOC | <u>DF</u> | Qualifiers |
| Acetone | <10 | 6.0 | 10 | 20 | 1.00 | U,ICH |
| Benzene | <0.50 | 0.14 | 0.50 | 1.0 | 1.00 | U |
| Bromodichloromethane | <0.50 | 0.21 | 0.50 | 5.0 | 1.00 | U |
| Bromoform | <1.0 | 0.50 | 1.0 | 10 | 1.00 | U |
| Bromomethane | <5.0 | 3.9 | 5.0 | 20 | 1.00 | U,IJ |
| 2-Butanone | <5.0 | 2.2 | 5.0 | 10 | 1.00 | U |
| Carbon Tetrachloride | <0.50 | 0.23 | 0.50 | 1.0 | 1.00 | U |
| Chlorobenzene | <0.50 | 0.17 | 0.50 | 5.0 | 1.00 | U |
| Chloroethane | <5.0 | 2.3 | 5.0 | 10 | 1.00 | U |
| Chloroform | <0.50 | 0.46 | 0.50 | 5.0 | 1.00 | U |
| Chloromethane | <2.0 | 1.8 | 2.0 | 10 | 1.00 | U |
| Dibromochloromethane | <0.50 | 0.25 | 0.50 | 1.0 | 1.00 | U |
| 1,2-Dibromo-3-Chloropropane | <2.0 | 1.2 | 2.0 | 10 | 1.00 | U |
| 1,2-Dibromoethane | <0.50 | 0.36 | 0.50 | 1.0 | 1.00 | U |
| 1,2-Dichlorobenzene | <0.50 | 0.46 | 0.50 | 1.0 | 1.00 | U |
| 1,3-Dichlorobenzene | <0.50 | 0.40 | 0.50 | 1.0 | 1.00 | U |
| 1,4-Dichlorobenzene | <0.50 | 0.43 | 0.50 | 1.0 | 1.00 | U |
| 1,1-Dichloroethane | <0.50 | 0.28 | 0.50 | 5.0 | 1.00 | U |
| 1,2-Dichloroethane | <0.50 | 0.24 | 0.50 | 1.0 | 1.00 | U |
| 1,1-Dichloroethene | <0.50 | 0.43 | 0.50 | 1.0 | 1.00 | U |
| c-1,2-Dichloroethene | <0.50 | 0.48 | 0.50 | 1.0 | 1.00 | U |
| t-1,2-Dichloroethene | <0.50 | 0.37 | 0.50 | 1.0 | 1.00 | U |
| 1,2-Dichloropropane | <0.50 | 0.42 | 0.50 | 5.0 | 1.00 | U |
| c-1,3-Dichloropropene | <0.50 | 0.25 | 0.50 | 1.0 | 1.00 | U |
| t-1,3-Dichloropropene | <0.50 | 0.25 | 0.50 | 1.0 | 1.00 | U |
| Ethylbenzene | <0.50 | 0.14 | 0.50 | 1.0 | 1.00 | U |
| Methylene Chloride | <1.0 | 0.64 | 1.0 | 5.0 | 1.00 | U |
| 4-Methyl-2-Pentanone | <5.0 | 4.4 | 5.0 | 10 | 1.00 | U |
| Styrene | <0.50 | 0.17 | 0.50 | 1.0 | 1.00 | U |
| 1,1,1,2-Tetrachloroethane | <0.50 | 0.40 | 0.50 | 1.0 | 1.00 | U |
| 1,1,2,2-Tetrachloroethane | <0.50 | 0.41 | 0.50 | 1.0 | 1.00 | U |
| Tetrachloroethene | <0.50 | 0.39 | 0.50 | 5.0 | 1.00 | U |
| Toluene | <0.50 | 0.24 | 0.50 | 1.0 | 1.00 | U |
| 1,2,4-Trichlorobenzene | <1.0 | 0.50 | 1.0 | 5.0 | 1.00 | U |
| 1,1,1-Trichloroethane | <0.50 | 0.30 | 0.50 | 5.0 | 1.00 | U |
| Hexachloro-1,3-Butadiene | <0.50 | 0.32 | 0.50 | 1.0 | 1.00 | U |



Environmental Science International, Inc. 354 Uluniu Street, Suite 304

Kailua, HI 96734-2500

1,4-Bromofluorobenzene

Date Received: Work Order:

Preparation:

Method: Units: 10/29/14 14-10-2234

EPA 5030C

GC/MS / EPA 8260B

ug/L

Project: Red Hill LTM 112066

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| *************************************** | | | | | | |
|---|---------------|----------------|------------|-----|-----------|------------|
| <u>Parameter</u> | <u>Result</u> | <u>DL</u> | LOD | LOQ | <u>DF</u> | Qualifiers |
| 1,1,2-Trichloroethane | <0.50 | 0.38 | 0.50 | 1.0 | 1.00 | U |
| Trichloroethene | <0.50 | 0.37 | 0.50 | 1.0 | 1.00 | U |
| 1,2,3-Trichloropropane | <1.0 | 0.64 | 1.0 | 5.0 | 1.00 | U |
| Vinyl Chloride | <0.50 | 0.30 | 0.50 | 1.0 | 1.00 | U |
| p/m-Xylene | <1.0 | 0.30 | 1.0 | 10 | 1.00 | U |
| o-Xylene | <0.50 | 0.23 | 0.50 | 1.0 | 1.00 | U |
| Methyl-t-Butyl Ether (MTBE) | <0.50 | 0.31 | 0.50 | 1.0 | 1.00 | U |
| Gasoline Range Organics | <30 | 26 | 30 | 50 | 1.00 | U |
| Surrogate | Rec. (%) | Control Limits | Qualifiers | | | |
| Dibromofluoromethane | 96 | 80-126 | | | | |
| 1,2-Dichloroethane-d4 | 100 | 80-134 | | | | |
| Toluene-d8 | 99 | 80-120 | | | | |
| Toluene-d8-TPPH | 96 | 88-112 | | | | |

80-120

97

10/29/14

14-10-2234

EPA 5030C



Analytical Report

Environmental Science International, Inc.

Date Received:

Work Order:

Kailua, HI 96734-2500 Preparation:

Method: GC/MS / EPA 8260B

Units: ug/L

Project: Red Hill LTM 112066 Page 7 of 8

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | | ate/Time QC Batch ID nalyzed |
|---|----------------------|------------------------|---------------|------------------|-----------------------|------------------------------|
| Method Blank | 099-13-057-72 | N/A | Aqueous | GC/MS OO | 10/31/14 10 | /31/14 141031L047 :31 |
| Comment(s): - Results were evaluated to | the MDL (DL), con | centrations >= t | o the MDL (DI | _) but < RL (LOC |), if found, are qual | ified with a "J" flag. |
| <u>Parameter</u> | <u>Result</u> | <u>DL</u> | LOD | LOC | <u>DF</u> | <u>Qualifiers</u> |
| Acetone | <10 | 6.0 | 10 | 20 | 1.00 | U |
| Benzene | <0.50 | 0.14 | 0.50 | 1.0 | 1.00 | U |
| Bromodichloromethane | <0.50 | 0.21 | 0.50 | 5.0 | 1.00 | U |
| Bromoform | <1.0 | 0.50 | 1.0 | 10 | 1.00 | U |
| Bromomethane | <5.0 | 3.9 | 5.0 | 20 | 1.00 | U |
| 2-Butanone | <5.0 | 2.2 | 5.0 | 10 | 1.00 | U |
| Carbon Tetrachloride | <0.50 | 0.23 | 0.50 | 1.0 | 1.00 | U |
| Chlorobenzene | <0.50 | 0.17 | 0.50 | 5.0 | 1.00 | U |
| Chloroethane | <5.0 | 2.3 | 5.0 | 10 | 1.00 | U |
| Chloroform | <0.50 | 0.46 | 0.50 | 5.0 | 1.00 | U |
| Chloromethane | <2.0 | 1.8 | 2.0 | 10 | 1.00 | U |
| Dibromochloromethane | <0.50 | 0.25 | 0.50 | 1.0 | 1.00 | U |
| 1,2-Dibromo-3-Chloropropane | <2.0 | 1.2 | 2.0 | 10 | 1.00 | U |
| 1,2-Dibromoethane | <0.50 | 0.36 | 0.50 | 1.0 | 1.00 | U |
| 1,2-Dichlorobenzene | <0.50 | 0.46 | 0.50 | 1.0 | 1.00 | U |
| 1,3-Dichlorobenzene | <0.50 | 0.40 | 0.50 | 1.0 | 1.00 | U |
| 1,4-Dichlorobenzene | <0.50 | 0.43 | 0.50 | 1.0 | 1.00 | U |
| 1,1-Dichloroethane | <0.50 | 0.28 | 0.50 | 5.0 | 1.00 | U |
| 1,2-Dichloroethane | <0.50 | 0.24 | 0.50 | 1.0 | 1.00 | U |
| 1,1-Dichloroethene | <0.50 | 0.43 | 0.50 | 1.0 | 1.00 | U |
| c-1,2-Dichloroethene | <0.50 | 0.48 | 0.50 | 1.0 | 1.00 | U |
| t-1,2-Dichloroethene | <0.50 | 0.37 | 0.50 | 1.0 | 1.00 | U |
| 1,2-Dichloropropane | <0.50 | 0.42 | 0.50 | 5.0 | 1.00 | U |
| c-1,3-Dichloropropene | <0.50 | 0.25 | 0.50 | 1.0 | 1.00 | U |
| t-1,3-Dichloropropene | <0.50 | 0.25 | 0.50 | 1.0 | 1.00 | U |
| Ethylbenzene | <0.50 | 0.14 | 0.50 | 1.0 | 1.00 | U |
| Methylene Chloride | <1.0 | 0.64 | 1.0 | 5.0 | 1.00 | U |
| 4-Methyl-2-Pentanone | <5.0 | 4.4 | 5.0 | 10 | 1.00 | U |
| Styrene | <0.50 | 0.17 | 0.50 | 1.0 | 1.00 | U |
| 1,1,1,2-Tetrachloroethane | <0.50 | 0.40 | 0.50 | 1.0 | 1.00 | U |
| 1,1,2,2-Tetrachloroethane | <0.50 | 0.41 | 0.50 | 1.0 | 1.00 | U |
| Tetrachloroethene | <0.50 | 0.39 | 0.50 | 5.0 | 1.00 | U |
| Toluene | <0.50 | 0.24 | 0.50 | 1.0 | 1.00 | U |
| 1,2,4-Trichlorobenzene | <1.0 | 0.50 | 1.0 | 5.0 | 1.00 | U |
| 1,1,1-Trichloroethane | <0.50 | 0.30 | 0.50 | 5.0 | 1.00 | U |
| Hexachloro-1,3-Butadiene | <0.50 | 0.32 | 0.50 | 1.0 | 1.00 | U |



Environmental Science International, Inc. 354 Uluniu Street, Suite 304

Kailua, HI 96734-2500

Date Received: Work Order:

Preparation:

Method: Units:

10/29/14

14-10-2234 EPA 5030C

GC/MS / EPA 8260B

ug/L

Project: Red Hill LTM 112066

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| <u>Parameter</u> | Result | <u>DL</u> | LOD | LOQ | DF | Qualifiers |
|-----------------------------|----------|----------------|------------|-----|------|------------|
| 1,1,2-Trichloroethane | <0.50 | 0.38 | 0.50 | 1.0 | 1.00 | U |
| Trichloroethene | <0.50 | 0.37 | 0.50 | 1.0 | 1.00 | U |
| 1,2,3-Trichloropropane | <1.0 | 0.64 | 1.0 | 5.0 | 1.00 | U |
| Vinyl Chloride | <0.50 | 0.30 | 0.50 | 1.0 | 1.00 | U |
| p/m-Xylene | <1.0 | 0.30 | 1.0 | 10 | 1.00 | U |
| o-Xylene | <0.50 | 0.23 | 0.50 | 1.0 | 1.00 | U |
| Methyl-t-Butyl Ether (MTBE) | <0.50 | 0.31 | 0.50 | 1.0 | 1.00 | U |
| Gasoline Range Organics | <30 | 26 | 30 | 50 | 1.00 | U |
| Surrogate | Rec. (%) | Control Limits | Qualifiers | | | |
| Dibromofluoromethane | 95 | 80-126 | | | | |
| 1,2-Dichloroethane-d4 | 99 | 80-134 | | | | |
| Toluene-d8 | 99 | 80-120 | | | | |
| Toluene-d8-TPPH | 95 | 88-112 | | | | |
| 1.4-Bromofluorobenzene | 96 | 80-120 | | | | |



Quality Control - Spike/Spike Duplicate

Environmental Science International, Inc.

Date Received:

Work Order:

10/29/14

Work Order:

14-10-2234

Kailua, HI 96734-2500

Preparation:

Method:

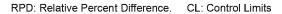
EPA 3005A Filt.

Method:

Project: Red Hill LTM 112066 Page 1 of 4

| Quality Control Sample ID | Туре | | Matrix | Ins | strument | Date Prepared | l Date Ana | lyzed | MS/MSD Bat | tch Number |
|---------------------------|-------------------------------|-----------------------|-------------|-------------|--------------|---------------|------------|-------|------------|------------|
| ES118 | Sample | | Aqueou | ıs ICI | P/MS 04 | 10/31/14 | 11/03/14 | 15:10 | 141031504 | |
| ES118 | Matrix Spike | | Aqueou | ıs ICI | P/MS 04 | 10/31/14 | 11/03/14 | 14:59 | 141031S04 | |
| ES118 | Matrix Spike I | Duplicate | Aqueou | is ICI | P/MS 04 | 10/31/14 | 11/03/14 | 15:01 | 141031504 | |
| Parameter | <u>Sample</u> <u>Conc.</u> | <u>Spike</u> Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Lead | ND | 100.0 | 109.0 | 109 | 109.2 | 109 | 80-120 | 0 | 0-20 | |

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Quality Control - Spike/Spike Duplicate

Environmental Science International, Inc.

Date Received:

Work Order:

10/29/14

Work Order:

Kailua, HI 96734-2500

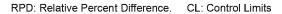
Preparation:

EPA 3510C

Method: EPA 8270C SIM PAHs

Project: Red Hill LTM 112066 Page 2 of 4

| Quality Control Sample ID | Туре | | Matrix | | Instrument | Date Prepared | l Date Ana | lyzed | MS/MSD Ba | tch Number |
|---------------------------|-------------------------------|------------------------------|-------------|------------|--------------|---------------|--------------------------|-------|-----------|------------|
| 14-10-2131-3 | Sample | | Aqueous | \$ | GC/MS AAA | 10/30/14 | 11/03/14 | 18:28 | 141030S08 | |
| 14-10-2131-3 | Matrix Spike | | Aqueous | | GC/MS AAA | 10/30/14 | 10/31/14 18:04 141030S08 | | | |
| 14-10-2131-3 | Matrix Spike | Duplicate | Aqueous | 5 | GC/MS AAA | 10/30/14 | 10/31/14 | 18:27 | 141030508 | |
| <u>Parameter</u> | <u>Sample</u> <u>Conc.</u> | <u>Spike</u> <u>Added</u> | MS Conc. | MS %Red | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| Naphthalene | 143.3 | 2.000 | 83.89 | 0 | 78.94 | 0 | 21-133 | 6 | 0-25 | 3 |
| 2-Methylnaphthalene | 42.74 | 2.000 | 34.71 | 0 | 32.01 | 0 | 21-140 | 8 | 0-25 | 3 |
| 1-Methylnaphthalene | 59.00 | 2.000 | 47.61 | 0 | 45.00 | 0 | 20-140 | 6 | 0-25 | 3 |
| Acenaphthylene | ND | 2.000 | 1.706 | 85 | 1.549 | 77 | 33-145 | 10 | 0-25 | |
| Acenaphthene | 0.5300 | 2.000 | 2.109 | 79 | 1.970 | 72 | 49-121 | 7 | 0-25 | |
| Fluorene | ND | 2.000 | 1.860 | 93 | 1.701 | 85 | 59-121 | 9 | 0-25 | |
| Phenanthrene | ND | 2.000 | 1.624 | 81 | 1.459 | 73 | 54-120 | 11 | 0-25 | |
| Anthracene | ND | 2.000 | 1.548 | 77 | 1.417 | 71 | 27-133 | 9 | 0-25 | |
| Fluoranthene | ND | 2.000 | 1.666 | 83 | 1.524 | 76 | 26-137 | 9 | 0-25 | |
| Pyrene | ND | 2.000 | 1.581 | 79 | 1.473 | 74 | 18-168 | 7 | 0-25 | |
| Benzo (a) Anthracene | ND | 2.000 | 1.551 | 78 | 1.422 | 71 | 33-143 | 9 | 0-25 | |
| Chrysene | ND | 2.000 | 1.540 | 77 | 1.402 | 70 | 17-168 | 9 | 0-25 | |
| Benzo (k) Fluoranthene | ND | 2.000 | 1.447 | 72 | 1.319 | 66 | 24-159 | 9 | 0-25 | |
| Benzo (b) Fluoranthene | ND | 2.000 | 1.372 | 69 | 1.248 | 62 | 24-159 | 9 | 0-25 | |
| Benzo (a) Pyrene | ND | 2.000 | 1.331 | 67 | 1.216 | 61 | 17-163 | 9 | 0-25 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 2.000 | 1.320 | 66 | 1.166 | 58 | 10-171 | 12 | 0-25 | |
| Dibenz (a,h) Anthracene | ND | 2.000 | 1.422 | 71 | 1.275 | 64 | 10-219 | 11 | 0-25 | |
| Benzo (g,h,i) Perylene | ND | 2.000 | 1.524 | 76 | 1.373 | 69 | 10-227 | 10 | 0-25 | |





Quality Control - Spike/Spike Duplicate

Environmental Science International, Inc.

Date Received: 10/29/14

354 Uluniu Street, Suite 304

Work Order: 14-10-2234

Kailua, HI 96734-2500

Preparation: EPA 5030C

Method: GC/MS / EPA 8260B

Project: Red Hill LTM 112066 Page 3 of 4

| 14-10-2387-2 Matrix Spite- pulseate Aqueous GC/MS OO 10/31/14 10/31/14 22:58 1410315031 + 1410315031 + 1410103712 14-10-2387-2 Matrix Spite- pulseate Aqueous GC/MS OO 10/31/14 10/31/14 22:58 1410315031 + 1010315031 + 1010315031 + 1010315031 + 1010315031 + 1010315031 + 1010315031 + 1010315031 + 1010315031 + 1010313031 + 101031 | Quality Control Sample ID | Туре | | Matrix | | Instrument | Date Prepared | d Date Analyzed MS/MSD Batch Number | | | | | |
|--|-----------------------------|--------------|-----------|-------------|-----------|-----------------|---------------|-------------------------------------|---------------------------|-----------|------------|--|--|
| 4-10-2367-2 Marity Spike burker Agenote Added Sociation Scheet Added MSD Scheet Added MSD Scheet Scheet Scheet Scheet Added MSD Scheet Scheet Added MSD Scheet Scheet Scheet Scheet Added MSD Scheet Scheet Added MSD Scheet Scheet Scheet Added MSD Scheet Scheet Scheet Added MSD Scheet Added< | 14-10-2387-2 | Sample | | Aqueou | S | GC/MS OO | 10/31/14 | 10/31/14 | 10/31/14 22:58 1410315031 | | | | |
| Parameter | 14-10-2387-2 | Matrix Spike | | Aqueou | S | GC/MS OO | 10/31/14 | 10/31/14 | 23:26 | 141031503 | 1 | | |
| Concord Conc | 14-10-2387-2 | Matrix Spike | Duplicate | Aqueou | s | GC/MS OO | 10/31/14 | 10/31/14 | 23:53 | 141031503 | 1 | | |
| Benzene ND 50.00 49.93 100 49.48 99 80-120 1 0-20 Bromodichloromethane ND 50.00 48.21 98 50.44 101 70-130 2 0-20 Bromodichloromethane ND 50.00 49.21 98 50.44 101 70-130 2 0-20 Bromofichane ND 50.00 59.57 119 59.62 119 30-145 17 0-20 2-Butanone ND 50.00 59.57 119 59.62 119 30-145 17 0-20 Chlorochane ND 50.00 48.15 96 48.88 98 60-125 1 0-20 Chlorochane ND 50.00 44.85 90 44.26 89 60-135 1 0-20 Chlorochane ND 50.00 49.33 99 50.66 101 60-135 1 0-20 Chloromethane ND 50.00 | <u>Parameter</u> | | | MS Conc. | MS %Re | MSD c. Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers | | |
| Bromodichloromethane ND 50.00 48.33 97 48.57 97 75.120 0 0.20 Bromoforform ND 50.00 49.21 88 50.44 101 70.130 2 0.20 Bromomethane ND 50.00 49.22 98 41.35 83 30.145 17 0.20 2-Butanone ND 50.00 51.03 102 49.80 100 65.140 2 0.20 Chlorobenzene ND 50.00 41.85 96 48.88 98 60.120 2 0.20 Chlorodethane ND 50.00 44.89 90 44.28 89 66.135 1 0.20 Chlorodethane ND 50.00 43.24 88 44.72 89 66.135 1 0.20 Chlorodethane ND 50.00 43.24 86 46.79 94 40.125 1 0.20 1,2-Dichromothare ND 50.00 | Acetone | ND | 50.00 | 67.72 | 135 | 76.03 | 152 | 40-140 | 12 | 0-20 | 3 | | |
| Bromotorm ND 50.00 49.21 98 50.44 101 70-130 2 0-20 Bromomethane ND 50.00 49.22 88 41.35 83 30-145 17 0-20 2-Butanone ND 50.00 59.57 119 59.62 119 30-150 0 0-20 Carbon Tetrachloride ND 50.00 48.15 96 48.88 98 80-120 2 0-20 Chlorobenzene ND 50.00 48.15 96 48.88 98 80-120 2 0-20 Chlorodethane ND 50.00 44.85 96 49.66 99 66-135 1 0-20 Chloromethane ND 50.00 44.19 88 44.72 89 40-125 1 0-20 Chloromethane ND 50.00 49.33 99 50.66 101 60-135 3 0-20 1,2-Dichrome-S-Chloropropane ND 5 | Benzene | ND | 50.00 | 49.93 | 100 | 49.48 | 99 | 80-120 | 1 | 0-20 | | | |
| Bromomethane ND 50.00 49.22 98 41.35 83 30-145 17 0-20 2-Butanone ND 50.00 59.57 119 59.62 119 30-150 0 0-20 Carbon Tetrachloride ND 50.00 51.03 102 49.80 100 65-140 2 0-20 Chloroethane ND 50.00 48.15 90 44.26 89 60-135 1 0-20 Chloroethane ND 50.00 43.30 90 44.26 89 65-135 1 0-20 Chloromethane ND 50.00 43.33 99 50.66 191 60-135 3 0-20 Dibromochloromethane ND 50.00 43.24 86 49.77 94 50-130 8 0-20 1,2-Dibromo-3-Chloropropane ND 50.00 48.24 86 49.97 94 50-130 4 0-20 1,2-Dibromo-3-Chloropropane <t< td=""><td>Bromodichloromethane</td><td>ND</td><td>50.00</td><td>48.33</td><td>97</td><td>48.57</td><td>97</td><td>75-120</td><td>0</td><td>0-20</td><td></td></t<> | Bromodichloromethane | ND | 50.00 | 48.33 | 97 | 48.57 | 97 | 75-120 | 0 | 0-20 | | | |
| 2-Butanone ND 50.00 59.57 119 59.62 119 30-150 0 0-20 Carbon Tetrachloride ND 50.00 \$1.03 102 49.80 100 65-140 2 0-20 Chlorobenzene ND 50.00 48.15 96 48.88 89 80-120 2 0-20 Chloromethane ND 50.00 50.30 101 49.66 89 65-135 1 0-20 Chloromethane ND 50.00 44.19 88 44.72 89 40-125 1 0-20 Chloromethane ND 50.00 49.33 99 50.66 101 60-135 3 0-20 1,2-Dibromo-3-Chloropropane ND 50.00 43.24 86 46.97 94 50-135 1 0-20 1,2-Dibrlomo-Bracene ND 50.00 47.25 95 48.04 96 75-125 0 0-20 1,3-Dichloroptenzene ND | Bromoform | ND | 50.00 | 49.21 | 98 | 50.44 | 101 | 70-130 | 2 | 0-20 | | | |
| Carbon Tetrachloride ND 50.00 51.03 102 49.80 100 65.140 2 0-20 Chlorobenzene ND 50.00 48.15 96 48.88 98 80-120 2 0-20 Chlorobentane ND 50.00 50.30 101 49.66 99 65-135 1 0-20 Chloromethane ND 50.00 44.19 88 44.72 89 40-125 1 0-20 Dibromochloromethane ND 50.00 49.33 99 50.66 101 60-135 3 0-20 1,2-Dibromo-3-Chloropropane ND 50.00 47.25 85 48.04 96 70-120 2 0-20 1,2-Dibriomoethane ND 50.00 47.25 85 48.04 96 75-125 0 0-20 1,2-Dibriomoethane ND 50.00 48.12 96 47.98 96 75-125 1 0-20 1,4-Dichlorobethane | Bromomethane | ND | 50.00 | 49.22 | 98 | 41.35 | 83 | 30-145 | 17 | 0-20 | | | |
| Chlorobenzene ND 50.00 48.15 96 48.88 98 80-120 2 0-20 Chlorobethane ND 50.00 44.85 90 44.26 89 60-135 1 0-20 Chloroform ND 50.00 50.30 101 49.66 99 65-135 1 0-20 Chloromethane ND 50.00 44.19 88 44.72 89 40-125 1 0-20 Dibromochloromethane ND 50.00 49.33 99 50.66 101 60-135 3 0-20 1,2-Dibromoethane ND 50.00 43.24 86 46.97 94 50-130 8 0-20 1,2-Dibromoethane ND 50.00 47.25 95 48.04 96 70-120 2 0-20 1,4-Dichlorobenzene ND 50.00 48.81 96 47.56 95 75-125 1 0-20 1,4-Dichlorobenzene ND | 2-Butanone | ND | 50.00 | 59.57 | 119 | 59.62 | 119 | 30-150 | 0 | 0-20 | | | |
| Chloroethane ND 50.00 44.85 90 44.26 89 60-135 1 0-20 Chloroform ND 50.00 50.30 101 49.66 99 65-135 1 0-20 Chloromethane ND 50.00 44.19 88 44.72 89 40-125 1 0-20 Dibromochloromethane ND 50.00 43.24 86 46.97 94 50-130 8 0-20 1,2-Dibromo-3-Chloropropane ND 50.00 47.25 95 48.04 96 70-120 2 0-20 1,2-Dibromoethane ND 50.00 47.25 95 48.04 96 70-120 2 0-20 1,2-Dichlorobenzene ND 50.00 48.17 96 47.86 96 75-125 0 0-20 1,4-Dichlorobenzene ND 50.00 48.86 94 47.56 95 75-125 1 0-20 1,1-Dichloroethane N | Carbon Tetrachloride | ND | 50.00 | 51.03 | 102 | 49.80 | 100 | 65-140 | 2 | 0-20 | | | |
| Chloroform ND 50.00 50.30 101 49.66 99 65-135 1 0-20 Chloromethane ND 50.00 44.19 88 44.72 89 40-125 1 0-20 Dibromochloromethane ND 50.00 49.33 99 50.66 101 60-135 3 0-20 1,2-Dibromo-3-Chloropropane ND 50.00 50.20 100 52.08 104 80-120 4 0-20 1,2-Dibriomoethane ND 50.00 47.25 95 48.04 96 70-120 2 0-20 1,3-Dichlorobenzene ND 50.00 48.17 96 47.98 96 75-125 0 0-20 1,4-Dichloroethane ND 50.00 48.12 96 47.56 95 75-125 1 0-20 1,1-Dichloroethane ND 50.00 48.89 94 46.82 94 70-130 0 0-20 1,1-Dichloroethane | Chlorobenzene | ND | 50.00 | 48.15 | 96 | 48.88 | 98 | 80-120 | 2 | 0-20 | | | |
| Chloromethane ND 50.00 44.19 88 44.72 89 40-125 1 0-20 Dibromochloromethane ND 50.00 49.33 99 50.66 101 60-135 3 0-20 1,2-Dibromochloromethane ND 50.00 43.24 86 46.97 94 50-130 8 0-20 1,2-Dibromochlane ND 50.00 47.25 95 48.04 96 70-120 2 0-20 1,3-Dichlorobenzene ND 50.00 47.25 95 48.04 96 70-120 2 0-20 1,4-Dichlorobenzene ND 50.00 48.817 96 47.50 96 75-125 0 0-20 1,4-Dichlorobenzene ND 50.00 48.86 94 47.50 95 75-125 1 0-20 1,1-Dichloroethane ND 50.00 48.89 94 46.82 94 70-130 1 0-20 1,1-Dichloroethane | Chloroethane | ND | 50.00 | 44.85 | 90 | 44.26 | 89 | 60-135 | 1 | 0-20 | | | |
| Dibromochloromethane ND 50.00 49.33 99 50.66 101 60-135 3 0-20 1,2-Dibromo-3-Chloropropane ND 50.00 43.24 86 46.97 94 50-130 8 0-20 1,2-Dibromoethane ND 50.00 50.20 100 52.08 104 80-120 4 0-20 1,2-Dibromoethane ND 50.00 47.25 95 48.04 96 70-120 2 0-20 1,3-Dichlorobenzene ND 50.00 48.17 96 47.98 96 75-125 0 0-20 1,4-Dichlorobenzene ND 50.00 48.17 96 47.50 95 75-125 1 0-20 1,4-Dichloroethane ND 50.00 48.12 96 47.56 95 75-125 1 0-20 1,1-Dichloroethane ND 50.00 48.82 96 47.56 95 75-125 1 0-20 1,2-Dichloroethane ND 50.00 48.89 94 46.82 94 70-130 1 0-20 1,1-Dichloroethane ND 50.00 51.13 102 50.25 101 70-130 1 0-20 1,2-Dichloroethene ND 50.00 51.13 102 50.25 101 70-125 2 0-20 1,2-Dichloroethene ND 50.00 51.62 103 48.81 96 75-125 0 0-20 1,2-Dichloroethene ND 50.00 51.62 103 52.11 104 70-130 1 0-20 1,2-Dichloroethene ND 50.00 51.62 103 52.11 104 70-130 1 0-20 1,2-Dichloroethene ND 50.00 51.62 103 52.11 104 70-130 1 0-20 1,2-Dichloroethene ND 50.00 51.62 103 52.11 104 70-130 1 0-20 1,1-Ja-Dichloropropene ND 50.00 48.66 99 49.72 99 75-125 0 0-20 1,1-Ja-Dichloropropene ND 50.00 47.52 95 49.08 98 55-140 3 0-20 2 Ethylbenzene ND 50.00 44.42 89 50.80 102 60-135 13 0-20 3 Ethylbenzene ND 50.00 48.42 89 50.80 102 60-135 13 0-20 3 Ethylbenzene ND 50.00 48.99 98 51.35 103 65-130 5 0-20 3 Ethylbenzene ND 50.00 48.99 98 51.35 103 65-130 5 0-20 4 Ethylbenzene ND 50.00 48.89 98 51.35 103 65-130 5 0-20 4 Ethylbenzene ND 50.00 48.89 98 51.35 103 65-130 5 0-20 5 Ethylbenzene ND 50.00 48.89 98 51.35 103 | Chloroform | ND | 50.00 | 50.30 | 101 | 49.66 | 99 | 65-135 | 1 | 0-20 | | | |
| 1,2-Dibromo-3-Chloropropane ND 50.00 43.24 86 46.97 94 50.130 8 0-20 1,2-Dibromoethane ND 50.00 50.20 100 52.08 104 80-120 4 0-20 1,2-Dichlorobenzene ND 50.00 47.25 95 48.04 96 75-125 0 0-20 1,4-Dichlorobenzene ND 50.00 48.17 96 47.98 96 75-125 1 0-20 1,4-Dichlorobenzene ND 50.00 48.12 96 47.56 95 75-125 1 0-20 1,4-Dichloroethane ND 50.00 48.81 96 47.56 95 70-135 1 0-20 1,2-Dichloroethane ND 50.00 48.85 100 50.61 101 70-130 0 0-20 -1,2-Dichloroethane ND 50.00 51.63 101 49.80 100 60-140 2 0-20 1,1-1,2-Dichl | Chloromethane | ND | 50.00 | 44.19 | 88 | 44.72 | 89 | 40-125 | 1 | 0-20 | | | |
| 1,2-Dibromoethane ND 50.00 50.20 100 52.08 104 80-120 4 0-20 1,2-Dichlorobenzene ND 50.00 47.25 95 48.04 96 70-120 2 0-20 1,3-Dichlorobenzene ND 50.00 48.17 96 47.98 96 75-125 0 0-20 1,4-Dichlorobenzene ND 50.00 48.86 94 47.50 95 75-125 1 0-20 1,1-Dichloroethane ND 50.00 48.12 96 47.56 95 70-135 1 0-20 1,2-Dichloroethane ND 50.00 48.85 100 50.61 101 70-130 1 0-20 -1,2-Dichloroethene ND 50.00 51.13 102 50.25 101 70-130 1 0-20 t-1,2-Dichloropropane ND 50.00 51.62 103 52.11 104 70-130 1 0-20 t-1,3-Dichloropro | Dibromochloromethane | ND | 50.00 | 49.33 | 99 | 50.66 | 101 | 60-135 | 3 | 0-20 | | | |
| 1,2-Dichlorobenzene ND 50.00 47.25 95 48.04 96 70-120 2 0-20 1,3-Dichlorobenzene ND 50.00 48.17 96 47.98 96 75-125 0 0-20 1,4-Dichlorobenzene ND 50.00 46.86 94 47.50 95 75-125 1 0-20 1,1-Dichloroethane ND 50.00 48.12 96 47.56 95 70-135 1 0-20 1,2-Dichloroethane ND 50.00 48.85 100 50.61 101 70-130 1 0-20 1,1-Dichloroethane ND 50.00 46.89 94 46.82 94 70-130 0 0-20 c-1,2-Dichloroethane ND 50.00 51.13 102 50.25 101 70-130 0 -20 t-1,2-Dichloroethane ND 50.00 48.06 96 48.21 96 75-125 0 0 -20 t-1 | 1,2-Dibromo-3-Chloropropane | ND | 50.00 | 43.24 | 86 | 46.97 | 94 | 50-130 | 8 | 0-20 | | | |
| 1,3-Dichlorobenzene ND 50.00 48.17 96 47.98 96 75-125 0 0-20 1,4-Dichlorobenzene ND 50.00 46.86 94 47.50 95 75-125 1 0-20 1,1-Dichloroethane ND 50.00 48.12 96 47.56 95 70-135 1 0-20 1,2-Dichloroethane ND 50.00 48.81 100 50.61 101 70-130 1 0-20 1,1-Dichloroethene ND 50.00 46.89 94 46.82 94 70-130 0 0-20 c-1,2-Dichloroethene ND 50.00 51.13 102 50.25 101 70-125 2 0-20 t-1,2-Dichloropropane ND 50.00 50.60 101 49.80 100 60-140 2 0-20 t-1,3-Dichloropropane ND 50.00 51.62 103 51.77 104 55-140 3 0-20 t-1,3-Dichloropropane ND 50.00 49.56 99 49.72 99 75-125 | 1,2-Dibromoethane | ND | 50.00 | 50.20 | 100 | 52.08 | 104 | 80-120 | 4 | 0-20 | | | |
| 1,4-Dichlorobenzene ND 50.00 46.86 94 47.50 95 75-125 1 0-20 1,1-Dichloroethane ND 50.00 48.12 96 47.56 95 70-135 1 0-20 1,2-Dichloroethane ND 50.00 49.85 100 50.61 101 70-130 1 0-20 1,1-Dichloroethene ND 50.00 46.89 94 46.82 94 70-130 0 0-20 c-1,2-Dichloroethene ND 50.00 51.13 102 50.25 101 70-125 2 0-20 t-1,2-Dichloropthene ND 50.00 50.60 101 49.80 100 60-140 2 0-20 t-2-Dichloroptropane ND 50.00 48.06 96 48.21 96 75-125 0 0-20 t-1,3-Dichloropropene ND 50.00 50.09 100 51.77 104 55-140 3 0-20 Ethylbenzene ND 50.00 47.52 95 49.08 98 55-140 <t< td=""><td>1,2-Dichlorobenzene</td><td>ND</td><td>50.00</td><td>47.25</td><td>95</td><td>48.04</td><td>96</td><td>70-120</td><td>2</td><td>0-20</td><td></td></t<> | 1,2-Dichlorobenzene | ND | 50.00 | 47.25 | 95 | 48.04 | 96 | 70-120 | 2 | 0-20 | | | |
| 1,1-Dichloroethane ND 50.00 48.12 96 47.56 95 70-135 1 0-20 1,2-Dichloroethane ND 50.00 49.85 100 50.61 101 70-130 1 0-20 1,1-Dichloroethene ND 50.00 46.89 94 46.82 94 70-130 0 0-20 c-1,2-Dichloroethene ND 50.00 51.13 102 50.25 101 70-125 2 0-20 t-1,2-Dichloroethene ND 50.00 50.60 101 49.80 100 60-140 2 0-20 t-1,2-Dichloroptopane ND 50.00 48.06 96 48.21 96 75-125 0 0-20 c-1,3-Dichloroptopene ND 50.00 51.62 103 52.11 104 70-130 1 0-20 Ethylbenzene ND 50.00 49.56 99 49.72 99 75-125 0 0-20 Methylere-Pentanone | 1,3-Dichlorobenzene | ND | 50.00 | 48.17 | 96 | 47.98 | 96 | 75-125 | 0 | 0-20 | | | |
| 1,2-Dichloroethane ND 50.00 49.85 100 50.61 101 70-130 1 0-20 1,1-Dichloroethene ND 50.00 46.89 94 46.82 94 70-130 0 0-20 c-1,2-Dichloroethene ND 50.00 51.13 102 50.25 101 70-125 2 0-20 t-1,2-Dichloroethene ND 50.00 50.60 101 49.80 100 60-140 2 0-20 t-2-Dichloropropane ND 50.00 48.06 96 48.21 96 75-125 0 0-20 c-1,3-Dichloropropane ND 50.00 51.62 103 52.11 104 70-130 1 0-20 t-1,3-Dichloropropene ND 50.00 50.09 100 51.77 104 55-140 3 0-20 Ethylbenzene ND 50.00 47.52 95 49.08 98 55-140 3 0-20 Methylene Chlori | 1,4-Dichlorobenzene | ND | 50.00 | 46.86 | 94 | 47.50 | 95 | 75-125 | 1 | 0-20 | | | |
| 1,1-Dichloroethene ND 50.00 46.89 94 46.82 94 70-130 0 0-20 c-1,2-Dichloroethene ND 50.00 51.13 102 50.25 101 70-125 2 0-20 t-1,2-Dichloroethene ND 50.00 50.60 101 49.80 100 60-140 2 0-20 1,2-Dichloropropane ND 50.00 48.06 96 48.21 96 75-125 0 0-20 c-1,3-Dichloropropene ND 50.00 51.62 103 52.11 104 70-130 1 0-20 t-1,3-Dichloropropene ND 50.00 50.09 100 51.77 104 55-140 3 0-20 Ethylbenzene ND 50.00 47.52 95 49.08 98 55-140 3 0-20 Methylene Chloride ND 50.00 44.42 89 50.80 102 60-135 13 0-20 Styrene ND 50.00 49.18 98 49.95 100 65-135 2 <td>1,1-Dichloroethane</td> <td>ND</td> <td>50.00</td> <td>48.12</td> <td>96</td> <td>47.56</td> <td>95</td> <td>70-135</td> <td>1</td> <td>0-20</td> <td></td> | 1,1-Dichloroethane | ND | 50.00 | 48.12 | 96 | 47.56 | 95 | 70-135 | 1 | 0-20 | | | |
| c-1,2-Dichloroethene ND 50.00 51.13 102 50.25 101 70-125 2 0-20 t-1,2-Dichloroethene ND 50.00 50.60 101 49.80 100 60-140 2 0-20 1,2-Dichloropropane ND 50.00 48.06 96 48.21 96 75-125 0 0-20 c-1,3-Dichloropropene ND 50.00 50.09 100 51.77 104 55-140 3 0-20 Ethylbenzene ND 50.00 49.56 99 49.72 99 75-125 0 0-20 Methylene Chloride ND 50.00 47.52 95 49.08 98 55-140 3 0-20 4-Methyl-2-Pentanone ND 50.00 44.42 89 50.80 102 60-135 13 0-20 Styrene ND 50.00 49.18 98 49.95 100 65-135 2 0-20 1,1,2,2-Tetrachloroethane <td>1,2-Dichloroethane</td> <td>ND</td> <td>50.00</td> <td>49.85</td> <td>100</td> <td>50.61</td> <td>101</td> <td>70-130</td> <td>1</td> <td>0-20</td> <td></td> | 1,2-Dichloroethane | ND | 50.00 | 49.85 | 100 | 50.61 | 101 | 70-130 | 1 | 0-20 | | | |
| t-1,2-Dichloroethene ND 50.00 50.60 101 49.80 100 60-140 2 0-20 1,2-Dichloropropane ND 50.00 48.06 96 48.21 96 75-125 0 0-20 c-1,3-Dichloropropene ND 50.00 51.62 103 52.11 104 70-130 1 0-20 t-1,3-Dichloropropene ND 50.00 50.09 100 51.77 104 55-140 3 0-20 Ethylbenzene ND 50.00 49.56 99 49.72 99 75-125 0 0-20 Methylene Chloride ND 50.00 47.52 95 49.08 98 55-140 3 0-20 4-Methyl-2-Pentanone ND 50.00 44.42 89 50.80 102 60-135 13 0-20 Styrene ND 50.00 49.18 98 49.95 100 65-135 2 0-20 1,1,2,2-Tetrachloroethane ND 50.00 50.43 101 50.94 102 80-130 <t< td=""><td>1,1-Dichloroethene</td><td>ND</td><td>50.00</td><td>46.89</td><td>94</td><td>46.82</td><td>94</td><td>70-130</td><td>0</td><td>0-20</td><td></td></t<> | 1,1-Dichloroethene | ND | 50.00 | 46.89 | 94 | 46.82 | 94 | 70-130 | 0 | 0-20 | | | |
| 1,2-Dichloropropane ND 50.00 48.06 96 48.21 96 75-125 0 0-20 c-1,3-Dichloropropene ND 50.00 51.62 103 52.11 104 70-130 1 0-20 t-1,3-Dichloropropene ND 50.00 50.09 100 51.77 104 55-140 3 0-20 Ethylbenzene ND 50.00 49.56 99 49.72 99 75-125 0 0-20 Methylene Chloride ND 50.00 47.52 95 49.08 98 55-140 3 0-20 4-Methyl-2-Pentanone ND 50.00 44.42 89 50.80 102 60-135 13 0-20 Styrene ND 50.00 49.18 98 49.95 100 65-135 2 0-20 1,1,2,2-Tetrachloroethane ND 50.00 50.43 101 50.94 102 80-130 1 0-20 Tetrachloroethane ND 50.00 50.71 101 52.32 105 45-150 3 | c-1,2-Dichloroethene | ND | 50.00 | 51.13 | 102 | 50.25 | 101 | 70-125 | 2 | 0-20 | | | |
| c-1,3-Dichloropropene ND 50.00 51.62 103 52.11 104 70-130 1 0-20 t-1,3-Dichloropropene ND 50.00 50.09 100 51.77 104 55-140 3 0-20 Ethylbenzene ND 50.00 49.56 99 49.72 99 75-125 0 0-20 Methylene Chloride ND 50.00 47.52 95 49.08 98 55-140 3 0-20 4-Methyl-2-Pentanone ND 50.00 44.42 89 50.80 102 60-135 13 0-20 Styrene ND 50.00 49.18 98 49.95 100 65-135 2 0-20 1,1,1,2-Tetrachloroethane ND 50.00 50.43 101 50.94 102 80-130 1 0-20 1,1,2,2-Tetrachloroethane ND 50.00 48.99 98 51.35 103 65-130 5 0-20 Tetrachloroethen | t-1,2-Dichloroethene | ND | 50.00 | 50.60 | 101 | 49.80 | 100 | 60-140 | 2 | 0-20 | | | |
| t-1,3-Dichloropropene ND 50.00 50.09 100 51.77 104 55-140 3 0-20 Ethylbenzene ND 50.00 49.56 99 49.72 99 75-125 0 0-20 Methylene Chloride ND 50.00 47.52 95 49.08 98 55-140 3 0-20 4-Methyl-2-Pentanone ND 50.00 44.42 89 50.80 102 60-135 13 0-20 Styrene ND 50.00 49.18 98 49.95 100 65-135 2 0-20 1,1,2-Tetrachloroethane ND 50.00 50.43 101 50.94 102 80-130 1 0-20 1,1,2,2-Tetrachloroethane ND 50.00 48.99 98 51.35 103 65-130 5 0-20 Tetrachloroethene ND 50.00 50.71 101 52.32 105 45-150 3 0-20 Toluene < | 1,2-Dichloropropane | ND | 50.00 | 48.06 | 96 | 48.21 | 96 | 75-125 | 0 | 0-20 | | | |
| Ethylbenzene ND 50.00 49.56 99 49.72 99 75-125 0 0-20 Methylene Chloride ND 50.00 47.52 95 49.08 98 55-140 3 0-20 4-Methyl-2-Pentanone ND 50.00 44.42 89 50.80 102 60-135 13 0-20 Styrene ND 50.00 49.18 98 49.95 100 65-135 2 0-20 1,1,2-Tetrachloroethane ND 50.00 50.43 101 50.94 102 80-130 1 0-20 1,1,2,2-Tetrachloroethane ND 50.00 48.99 98 51.35 103 65-130 5 0-20 Tetrachloroethene ND 50.00 50.71 101 52.32 105 45-150 3 0-20 Tolluene ND 50.00 48.58 97 49.23 98 75-120 1 0-20 | c-1,3-Dichloropropene | ND | 50.00 | 51.62 | 103 | 52.11 | 104 | 70-130 | 1 | 0-20 | | | |
| Methylene Chloride ND 50.00 47.52 95 49.08 98 55-140 3 0-20 4-Methyl-2-Pentanone ND 50.00 44.42 89 50.80 102 60-135 13 0-20 Styrene ND 50.00 49.18 98 49.95 100 65-135 2 0-20 1,1,2-Tetrachloroethane ND 50.00 50.43 101 50.94 102 80-130 1 0-20 1,1,2,2-Tetrachloroethane ND 50.00 48.99 98 51.35 103 65-130 5 0-20 Tetrachloroethene ND 50.00 50.71 101 52.32 105 45-150 3 0-20 Toluene ND 50.00 48.58 97 49.23 98 75-120 1 0-20 | t-1,3-Dichloropropene | ND | 50.00 | 50.09 | 100 | 51.77 | 104 | 55-140 | 3 | 0-20 | | | |
| 4-Methyl-2-Pentanone ND 50.00 44.42 89 50.80 102 60-135 13 0-20 Styrene ND 50.00 49.18 98 49.95 100 65-135 2 0-20 1,1,1,2-Tetrachloroethane ND 50.00 50.43 101 50.94 102 80-130 1 0-20 1,1,2,2-Tetrachloroethane ND 50.00 48.99 98 51.35 103 65-130 5 0-20 Tetrachloroethene ND 50.00 50.71 101 52.32 105 45-150 3 0-20 Tolluene ND 50.00 48.58 97 49.23 98 75-120 1 0-20 | Ethylbenzene | ND | 50.00 | 49.56 | 99 | 49.72 | 99 | 75-125 | 0 | 0-20 | | | |
| Styrene ND 50.00 49.18 98 49.95 100 65-135 2 0-20 1,1,1,2-Tetrachloroethane ND 50.00 50.43 101 50.94 102 80-130 1 0-20 1,1,2,2-Tetrachloroethane ND 50.00 48.99 98 51.35 103 65-130 5 0-20 Tetrachloroethene ND 50.00 50.71 101 52.32 105 45-150 3 0-20 Toluene ND 50.00 48.58 97 49.23 98 75-120 1 0-20 | Methylene Chloride | ND | 50.00 | 47.52 | 95 | 49.08 | 98 | 55-140 | 3 | 0-20 | | | |
| 1,1,1,2-Tetrachloroethane ND 50.00 50.43 101 50.94 102 80-130 1 0-20 1,1,2,2-Tetrachloroethane ND 50.00 48.99 98 51.35 103 65-130 5 0-20 Tetrachloroethene ND 50.00 50.71 101 52.32 105 45-150 3 0-20 Toluene ND 50.00 48.58 97 49.23 98 75-120 1 0-20 | 4-Methyl-2-Pentanone | ND | 50.00 | 44.42 | 89 | 50.80 | 102 | 60-135 | 13 | 0-20 | | | |
| 1,1,2,2-Tetrachloroethane ND 50.00 48.99 98 51.35 103 65-130 5 0-20 Tetrachloroethene ND 50.00 50.71 101 52.32 105 45-150 3 0-20 Toluene ND 50.00 48.58 97 49.23 98 75-120 1 0-20 | Styrene | ND | 50.00 | 49.18 | 98 | 49.95 | 100 | 65-135 | 2 | 0-20 | | | |
| Tetrachloroethene ND 50.00 50.71 101 52.32 105 45-150 3 0-20 Toluene ND 50.00 48.58 97 49.23 98 75-120 1 0-20 | 1,1,1,2-Tetrachloroethane | ND | 50.00 | 50.43 | 101 | 50.94 | 102 | 80-130 | 1 | 0-20 | | | |
| Toluene ND 50.00 48.58 97 49.23 98 75-120 1 0-20 | 1,1,2,2-Tetrachloroethane | ND | 50.00 | 48.99 | 98 | 51.35 | 103 | 65-130 | 5 | 0-20 | | | |
| | Tetrachloroethene | ND | 50.00 | 50.71 | 101 | 52.32 | 105 | 45-150 | 3 | 0-20 | | | |
| 1,2,4-Trichlorobenzene ND 50.00 44.33 89 46.15 92 65-135 4 0-20 | Toluene | ND | 50.00 | 48.58 | 97 | 49.23 | 98 | 75-120 | 1 | 0-20 | | | |
| | 1,2,4-Trichlorobenzene | ND | 50.00 | 44.33 | 89 | 46.15 | 92 | 65-135 | 4 | 0-20 | | | |
| 1,1,1-Trichloroethane ND 50.00 50.20 100 48.87 98 65-130 3 0-20 | 1,1,1-Trichloroethane | ND | 50.00 | 50.20 | 100 | 48.87 | 98 | 65-130 | 3 | 0-20 | | | |



Quality Control - Spike/Spike Duplicate

Environmental Science International, Inc.

Date Received:

Work Order:

14-10-2234

Kailua, HI 96734-2500

Preparation:

Method:

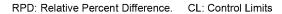
Date Received:

10/29/14

EPA 5030C

Project: Red Hill LTM 112066 Page 4 of 4

| <u>Parameter</u> | Sample Conc. | <u>Spike</u> Added | MS Conc. | MS %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
|-----------------------------|-----------------|-----------------------|-------------|-------------|--------------|--------------|----------|-----|--------|------------|
| Hexachloro-1,3-Butadiene | ND | 50.00 | 45.69 | 91 | 45.79 | 92 | 50-140 | 0 | 0-20 | |
| 1,1,2-Trichloroethane | ND | 50.00 | 49.59 | 99 | 51.28 | 103 | 75-125 | 3 | 0-20 | |
| Trichloroethene | ND | 50.00 | 47.79 | 96 | 47.66 | 95 | 70-125 | 0 | 0-20 | |
| 1,2,3-Trichloropropane | ND | 50.00 | 47.54 | 95 | 51.06 | 102 | 75-125 | 7 | 0-20 | |
| Vinyl Chloride | ND | 50.00 | 47.42 | 95 | 46.59 | 93 | 50-145 | 2 | 0-20 | |
| p/m-Xylene | ND | 100.0 | 98.13 | 98 | 98.70 | 99 | 75-130 | 1 | 0-20 | |
| o-Xylene | ND | 50.00 | 49.82 | 100 | 49.94 | 100 | 80-120 | 0 | 0-20 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 50.00 | 47.63 | 95 | 49.54 | 99 | 65-125 | 4 | 0-20 | |





Quality Control - PDS

Environmental Science International, Inc.

Date Received:

Work Order:

14-10-2234

Kailua, HI 96734-2500

Preparation:

Method:

EPA 3005A Filt.

Method:

Project: Red Hill LTM 112066 Page 1 of 1

| Quality Control Sample ID | Туре | M | 1atrix | Instrument | Date Prepared I | Date Analyzed | PDS/PDSD Batch Number | |
|---------------------------|--------|--------------|-------------|------------|-----------------|----------------|--------------------------|--|
| ES118 | Sample | А | queous | ICP/MS 04 | 10/31/14 00:00 | 11/03/14 15:10 | 141031504 | |
| ES118 | PDS | Α | queous | ICP/MS 04 | 10/31/14 00:00 | 11/03/14 15:03 | 141031S04 | |
| <u>Parameter</u> | | Sample Conc. | Spike Added | PDS Conc. | . PDS %Red | <u>%Rec. C</u> | <u>Qualifiers</u> | |
| Lead | | ND | 100.0 | 102.6 | 103 | 75-125 | | |



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Page 1 of 5

0-11



Project: Red Hill LTM 112066

TPH as Diesel

2000

2296

Quality Control - LCS/LCSD

Environmental Science International, Inc.

Date Received:

Work Order:

14-10-2234

Kailua, HI 96734-2500

Preparation:

Method:

EPA 8015B (M)

| Quality Control Sample ID | Туре | Matrix | Instrument | Date Prepared | Date Analyzed | LCS/LCSD B | atch Number |
|---------------------------|---------------|----------------------------|------------|-------------------|----------------|------------|-------------------|
| 099-15-516-206 | LCS | Aqueous | GC 45 | 10/30/14 | 10/31/14 07:20 | 141030B06 | |
| 099-15-516-206 | LCSD | Aqueou: | GC 45 | 10/30/14 | 10/31/14 07:37 | 141030B06 | |
| <u>Parameter</u> | Spike Added L | CS Conc. LCS <u>%</u> R | | LCSD %Re %Rec. | ec. CL RPD | RPD CL | <u>Qualifiers</u> |

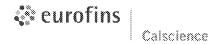
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111

60-132

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115



Project: Red Hill LTM 112066

Quality Control - LCS

Environmental Science International, Inc.

Date Received:

Work Order:

14-10-2234

Kailua, HI 96734-2500

Preparation:

Method:

EPA 3005A Filt.

Method:

Page 2 of 5

| Quality Control Sample ID | Туре | Matrix | Instrument [| Date Prepared | Date Analyzed LC | S Batch Number |
|---------------------------|------|-------------|-----------------|---------------|--------------------|-------------------|
| 099-14-497-104 | LCS | Aqueous | ICP/MS 04 1 | 0/31/14 | 11/03/14 14:50 14: | 1031L04F |
| <u>Parameter</u> | | Spike Added | Conc. Recovered | d LCS %Re | ec. %Rec. CL | <u>Qualifiers</u> |
| Lead | | 100.0 | 102.6 | 103 | 80-120 | |



Quality Control - LCS

Environmental Science International, Inc.

354 Uluniu Street, Suite 304 Kailua, HI 96734-2500 Date Received:

10/29/14 14-10-2234

Work Order: Preparation:

EPA 3510C

Method:

EPA 8270C SIM PAHs

| Project: Red Hill LTM 11 | 2066 | | | | | Page 3 of 5 | |
|---------------------------|------|---------|------------|---------------|----------------|------------------|--|
| Quality Control Sample ID | Туре | Matrix | Instrument | Date Prepared | Date Analyzed | LCS Batch Number | |
| 099-15-148-67 | LCS | Aqueous | GC/MS AAA | 10/30/14 | 10/31/14 17:40 | 141030L08 | |

| addity Control Cample 15 | 1 9 PC | Many | modunich | Date i repared Date | Analyzed Loo ba | ttori ivarribei |
|---------------------------|--------|-------------|-----------------|---------------------|-------------------|-------------------|
| 099-15-148-67 | LCS | Aqueous | GC/MS AAA | 10/30/14 10/3 | 1/14 17:40 141030 | L08 |
| <u>Parameter</u> | | Spike Added | Conc. Recovered | ed LCS %Rec. | %Rec. CL | <u>Qualifiers</u> |
| Naphthalene | | 2.000 | 1.394 | 70 | 21-133 | |
| 2-Methylnaphthalene | | 2.000 | 1.382 | 69 | 21-140 | |
| 1-Methylnaphthalene | | 2.000 | 1.409 | 70 | 20-140 | |
| Acenaphthylene | | 2.000 | 1.394 | 70 | 33-145 | |
| Acenaphthene | | 2.000 | 1.451 | 73 | 55-121 | |
| Fluorene | | 2.000 | 1.539 | 77 | 59-121 | |
| Phenanthrene | | 2.000 | 1.447 | 72 | 54-120 | |
| Anthracene | | 2.000 | 1.308 | 65 | 27-133 | |
| Fluoranthene | | 2.000 | 1.538 | 77 | 26-137 | |
| Pyrene | | 2.000 | 1.480 | 74 | 45-129 | |
| Benzo (a) Anthracene | | 2.000 | 1.398 | 70 | 33-143 | |
| Chrysene | | 2.000 | 1.425 | 71 | 17-168 | |
| Benzo (k) Fluoranthene | | 2.000 | 1.412 | 71 | 24-159 | |
| Benzo (b) Fluoranthene | | 2.000 | 1.322 | 66 | 24-159 | |
| Benzo (a) Pyrene | | 2.000 | 1.150 | 58 | 17-163 | |
| Indeno (1,2,3-c,d) Pyrene | | 2.000 | 1.251 | 63 | 25-175 | |
| Dibenz (a,h) Anthracene | | 2.000 | 1.391 | 70 | 25-175 | |
| Benzo (g,h,i) Perylene | | 2.000 | 1.514 | 76 | 25-157 | |



Quality Control - LCS/LCSD

Environmental Science International, Inc.

Date Received:

Work Order:

10/29/14

Work Order:

Kailua, HI 96734-2500

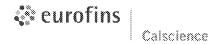
Preparation:

EPA 5030C

Method: GC/MS / EPA 8260B

Project: Red Hill LTM 112066 Page 4 of 5

| Quality Control Sample ID Type | | Mat | rix | Instrument | Date Pre | Date Prepared Date Analyzed LCS/LCSD Batch Nun | | | |
|--------------------------------|-------------|-----------|--------------|------------|---------------|--|------------|------------|------------|
| 099-13-057-72 | LCS | Aqı | ieous | GC/MS OO | 10/31/14 | 10/3 | 1/14 21:09 | 141031L047 | |
| 099-13-057-72 | LCSD | Aqı | ieous | GC/MS OO | 10/31/14 | 10/3 | 1/14 21:37 | 141031L047 | |
| <u>Parameter</u> | Spike Added | LCS Conc. | LCS %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | <u>RPD</u> | RPD CL | Qualifiers |
| Acetone | 50.00 | 66.09 | 132 | N/A | N/A | 40-140 | N/A | 0-20 | |
| Benzene | 50.00 | 50.90 | 102 | N/A | N/A | 80-120 | N/A | 0-20 | |
| Bromodichloromethane | 50.00 | 49.29 | 99 | N/A | N/A | 75-120 | N/A | 0-20 | |
| Bromoform | 50.00 | 50.67 | 101 | N/A | N/A | 70-130 | N/A | 0-20 | |
| Bromomethane | 50.00 | 43.60 | 87 | N/A | N/A | 30-145 | N/A | 0-20 | |
| 2-Butanone | 50.00 | 55.76 | 112 | N/A | N/A | 30-150 | N/A | 0-20 | |
| Carbon Tetrachloride | 50.00 | 52.82 | 106 | N/A | N/A | 65-140 | N/A | 0-20 | |
| Chlorobenzene | 50.00 | 50.31 | 101 | N/A | N/A | 80-120 | N/A | 0-20 | |
| Chloroethane | 50.00 | 46.45 | 93 | N/A | N/A | 60-135 | N/A | 0-20 | |
| Chloroform | 50.00 | 51.26 | 103 | N/A | N/A | 65-135 | N/A | 0-20 | |
| Chloromethane | 50.00 | 46.27 | 93 | N/A | N/A | 40-125 | N/A | 0-20 | |
| Dibromochloromethane | 50.00 | 50.80 | 102 | N/A | N/A | 60-135 | N/A | 0-20 | |
| 1,2-Dibromo-3-Chloropropane | 50.00 | 47.64 | 95 | N/A | N/A | 50-130 | N/A | 0-20 | |
| 1,2-Dibromoethane | 50.00 | 52.19 | 104 | N/A | N/A | 80-120 | N/A | 0-20 | |
| 1,2-Dichlorobenzene | 50.00 | 49.11 | 98 | N/A | N/A | 70-120 | N/A | 0-20 | |
| 1,3-Dichlorobenzene | 50.00 | 49.80 | 100 | N/A | N/A | 75-125 | N/A | 0-20 | |
| 1,4-Dichlorobenzene | 50.00 | 49.36 | 99 | N/A | N/A | 75-125 | N/A | 0-20 | |
| 1,1-Dichloroethane | 50.00 | 50.29 | 101 | N/A | N/A | 70-135 | N/A | 0-20 | |
| 1,2-Dichloroethane | 50.00 | 51.26 | 103 | N/A | N/A | 70-130 | N/A | 0-20 | |
| 1,1-Dichloroethene | 50.00 | 50.07 | 100 | N/A | N/A | 70-130 | N/A | 0-20 | |
| c-1,2-Dichloroethene | 50.00 | 52.18 | 104 | N/A | N/A | 70-125 | N/A | 0-20 | |
| t-1,2-Dichloroethene | 50.00 | 53.26 | 107 | N/A | N/A | 60-140 | N/A | 0-20 | |
| 1,2-Dichloropropane | 50.00 | 49.39 | 99 | N/A | N/A | 75-125 | N/A | 0-20 | |
| c-1,3-Dichloropropene | 50.00 | 53.06 | 106 | N/A | N/A | 70-130 | N/A | 0-20 | |
| t-1,3-Dichloropropene | 50.00 | 53.91 | 108 | N/A | N/A | 55-140 | N/A | 0-20 | |
| Ethylbenzene | 50.00 | 52.31 | 105 | N/A | N/A | 75-125 | N/A | 0-20 | |
| Methylene Chloride | 50.00 | 50.23 | 100 | N/A | N/A | 55-140 | N/A | 0-20 | |
| 4-Methyl-2-Pentanone | 50.00 | 49.85 | 100 | N/A | N/A | 60-135 | N/A | 0-20 | |
| Styrene | 50.00 | 51.56 | 103 | N/A | N/A | 65-135 | N/A | 0-20 | |
| 1,1,1,2-Tetrachloroethane | 50.00 | 51.98 | 104 | N/A | N/A | 80-130 | N/A | 0-20 | |
| 1,1,2,2-Tetrachloroethane | 50.00 | 50.95 | 102 | N/A | N/A | 65-130 | N/A | 0-20 | |
| Tetrachloroethene | 50.00 | 51.34 | 103 | N/A | N/A | 45-150 | N/A | 0-20 | |
| Toluene | 50.00 | 50.99 | 102 | N/A | N/A | 75-120 | N/A | 0-20 | |
| 1,2,4-Trichlorobenzene | 50.00 | 47.93 | 96 | N/A | N/A | 65-135 | N/A | 0-20 | |
| 1,1,1-Trichloroethane | 50.00 | 52.27 | 105 | N/A | N/A | 65-130 | N/A | 0-20 | |
| Hexachloro-1,3-Butadiene | 50.00 | 49.37 | 99 | N/A | N/A | 50-140 | N/A | 0-20 | |



Quality Control - LCS/LCSD

Environmental Science International, Inc.

Date Received:

Work Order:

14-10-2234

Kailua, HI 96734-2500

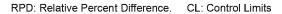
Preparation:

Method:

GC/MS / EPA 8260B

Project: Red Hill LTM 112066 Page 5 of 5

| <u>Parameter</u> | Spike Added | LCS Conc. | <u>LCS</u> %Rec. | LCSD Conc. | LCSD %Rec. | %Rec. CL | RPD | RPD CL | <u>Qualifiers</u> |
|-----------------------------|-------------|-----------|---------------------|------------|---------------|----------|-----|--------|-------------------|
| 1,1,2-Trichloroethane | 50.00 | 50.61 | 101 | N/A | N/A | 75-125 | N/A | 0-20 | |
| Trichloroethene | 50.00 | 50.27 | 101 | N/A | N/A | 70-125 | N/A | 0-20 | |
| 1,2,3-Trichloropropane | 50.00 | 50.55 | 101 | N/A | N/A | 75-125 | N/A | 0-20 | |
| Vinyl Chloride | 50.00 | 49.26 | 99 | N/A | N/A | 50-145 | N/A | 0-20 | |
| p/m-Xylene | 100.0 | 103.2 | 103 | N/A | N/A | 75-130 | N/A | 0-20 | |
| o-Xylene | 50.00 | 51.96 | 104 | N/A | N/A | 80-120 | N/A | 0-20 | |
| Methyl-t-Butyl Ether (MTBE) | 50.00 | 50.09 | 100 | N/A | N/A | 65-125 | N/A | 0-20 | |
| Gasoline Range Organics | 1000 | 1009 | 101 | 1025 | 103 | 80-120 | 2 | 0-20 | |





Sample Analysis Summary Report

| Work Order: 14-10-2234 | | | | Page 1 of 1 |
|------------------------|-----------------|------------|------------|---------------------|
| <u>Method</u> | Extraction | Chemist ID | Instrument | Analytical Location |
| EPA 6020 | EPA 3005A Filt. | 598 | ICP/MS 04 | 1 |
| EPA 8015B (M) | EPA 3510C | 682 | GC 45 | 1 |
| EPA 8270C SIM PAHs | EPA 3510C | 923 | GC/MS AAA | 1 |
| GC/MS / EPA 8260B | EPA 5030C | 849 | GC/MS OO | 2 |

Location 1: 7440 Lincoln Way, Garden Grove, CA 92841 Location 2: 7445 Lampson Avenue, Garden Grove, CA 92841





Glossary of Terms and Qualifiers

Work Order: 14-10-2234 Page 1 of 1

| Qualifiers | <u>Definition</u> |
|------------|--|
| * | See applicable analysis comment. |
| < | Less than the indicated value. |
| > | Greater than the indicated value. |
| 1 | Surrogate compound recovery was out of control due to a required sample dilution. Therefore, the sample data was reported without further clarification. |
| 2 | Surrogate compound recovery was out of control due to matrix interference. The associated method blank surrogate spike compound was in control and, therefore, the sample data was reported without further clarification. |
| 3 | Recovery of the Matrix Spike (MS) or Matrix Spike Duplicate (MSD) compound was out of control due to suspected matrix interference. The associated LCS recovery was in control. |
| 4 | The MS/MSD RPD was out of control due to suspected matrix interference. |
| 5 | The PDS/PDSD or PES/PESD associated with this batch of samples was out of control due to suspected matrix interference. |
| 6 | Surrogate recovery below the acceptance limit. |
| 7 | Surrogate recovery above the acceptance limit. |
| В | Analyte was present in the associated method blank. |
| BU | Sample analyzed after holding time expired. |
| BV | Sample received after holding time expired. |
| DL | The Detection Limit (DL) is the smallest analyte concentration that can be demonstrated to be different from zero or a blank concentration at the 99% level of confidence. |
| E | Concentration exceeds the calibration range. |
| ET | Sample was extracted past end of recommended max. holding time. |
| HD | The chromatographic pattern was inconsistent with the profile of the reference fuel standard. |
| ICH | Initial calibration verification recovery is above the control limit for this analyte. |
| ICJ | Initial calibration verification recovery is below the control limit for this analyte. |
| IH | Calibration verification recovery is above the control limit for this analyte. |
| IJ | Calibration verification recovery is below the control limit for this analyte. |
| J | Analyte was detected at a concentration below the LOQ and above the DL. Reported value is estimated. |
| JA | Analyte positively identified but quantitation is an estimate. |
| LOD | The Limit of Detection (LOD) is the smallest amount or concentration of a substance that must be present in a sample in order to be detected at 99% confidence level. |
| LOQ | The Limit of Quantitation (LOQ) is the lowest concentration of a substance that produces a quantitative result within specified limits of precision and bias. |
| Q | Spike recovery and RPD control limits do not apply resulting from the parameter concentration in the sample exceeding the spike concentration by a factor of four or greater. |
| SG | The sample extract was subjected to Silica Gel treatment prior to analysis. |
| U | Undetected at Detection Limit (DL) and is reported as less than the Limit of Detection (LOD). |
| Χ | % Recovery and/or RPD out-of-range. |
| Z | Analyte presence was not confirmed by second column or GC/MS analysis. |
| | Solid - Unless otherwise indicated, solid sample data is reported on a wet weight basis, not corrected for % moisture. All OC results are |

Solid - Unless otherwise indicated, solid sample data is reported on a wet weight basis, not corrected for % moisture. All QC results are reported on a wet weight basis.

Any parameter identified in 40CFR Part 136.3 Table II that is designated as "analyze immediately" with a holding time of <= 15 minutes (40CFR-136.3 Table II, footnote 4), is considered a "field" test and the reported results will be qualified as being received outside of the stated holding time unless received at the laboratory within 15 minutes of the collection time.

A calculated total result (Example: Total Pesticides) is the summation of each component concentration and/or, if "J" flags are reported, estimated concentration. Component concentrations showing not detected (ND) are summed into the calculated total result as zero concentrations.

| , 1 | | | relation entres | | | | | | | | | | unomicoran promissiones | | Page | 29 (| of 31 |
|--|-------------------------------|------------------------------------|---|--|--|--|--|-----------------------------|--------------------|--|--|----------|-------------------------|--|--|--------------------------------------|--------------------------------------|
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| 128/14 | P.O. NO.: | SAMPLER(S): (PRINT) | 以方方 | ANALYSES | as needed. | | MIS 0728 18 □ X747\01 18.812 □ 9917 | X 07S8 | steM SST | - | X | X | | | Date: Vofosji S | Date: | Date; |
| Date O | ~ ~ ~ | 5 9 9 | | | check box or fill in blank as ne | e Core | rneT 🗆 encO r | 35) 🗆 Er | Sygens (50: | | | | | | and the control of th | | Salkt Salkt |
| 0#/LAB USE ONLY [4-10-2234 | CLIENT PROJECT NAME / NUMBER | ROJECT CONTACT: | は中である | | Please check box | | 85e0 🗆 🗆 Ce-C44 | □ 38TM | TPH | | 8 | X | | | contrare l'acception annoise descriptions de construction de la contrare de la contrare de la contrare de la c | | |
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| -5494 s Christi | | | 734 ZIP: | Perienes | ARD | LOG CODE | | | Unprese | Sectional Contraction Contract | У. | × | | | Received by: (Sig | Received by: (Sig | Received by: (Sig |
| 7440 Lincoln Way, Garden Grove, CA 92841-1427 • (714) 895-5494 Other locations: Concord, San Luis Obispo, Houston, and Corpus Christi For courier service / sample drop off information, contact sales@calscience.com or call us. | | | STATE: 96 | EMALE MERCELOPERCE, COM, APPERCEPTIONES COM | :D"): 'S (XSTANDARD | Available of the Control of the Cont | を呼 | | MATRIX OF | 200 | ************************************** | 1 1 | | | | | |
| Nay, Garden Grove, CA 92841-1427 • (7 Concord, San Luis Obispo, Houston, and courier service / sample drop off informati contact sales@calscience.com or call us. | | まり30年 | Ţ | Occience. | AT not "STANDARD" HR | eraktiona | ing anal | | SAMPLING E TIME | 0011 141/ | 1000 FE | 12/30 | | | | | |
| oln Way, Garden Grove, CA 92841-1427 • (714) ons: Concord, San Luis Obispo, Houston, and Cc For courier service / sample drop off information, contact sales@calscience.com or call us. | | え、あ | ancipientamannes earlinessessiffficturents ancientosite properties ancientosite ancientos ancientos ancientos a | Tations C | ss may apply to any TAT | on a state and state of state | monitar | | S | h/82/01 | | 10/26/14 | | | | | |
| 7440 Lincoln Other locations | LABORATORY CLIENT: ESI | ADDRESS: 354 Ulyning St., Suit 304 | OITY: Kailna | TEL (808) 261-0740 | TURNAROUND TIME (Rush surcharges may apply to any TAT not "STANDARD"): SAME DAY 124 HR 5 DAYS | COELT EDF GLOBAL ID | special instructions: Degular (full) fim monitaring analytelist | | SAMPLEID | 65TRIP | かいこ | 155 - SX | | | Relinquished by: (Signature) | Relinquished by: (Signature) | Relinquished by: (Signature) |
| | BORA | ODRES | IIX: | | URNAF]SAME | COE | PECIAL PECIAL | | LAB USE | 100 | N | ~ | | | Selinqu Selinqu | Selinqu | elinqu |

ORIGIN ID:HNLA

SHIP DATE: 280CT14 ACTWGT: 52.5 LB CAD: /POS1525 DIMS: 26x15x14 IN

BILL RECIPIENT

UNITED STATES US

TO

CALSCIENCE 7440 LINCOLN WAY

GARDEN GROVE CA 92841

(714) 895 - 5494 INU: PO:

REF:

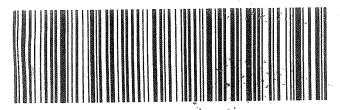


TRK# 8045 5791 7375

WED - 29 OCT AA STANDARD OVERNIGHT

WZ APVA

92841 ca=us SNA



(223^H)

Cooler <u>/</u> of <u>/</u>



Calscience

WORK ORDER #: 14-10-2 2 3 任

SAMPLE RECEIPT FORM

| CLIENT: ESI | DATE: _ | 10/29 | /14 |
|---|--------------|----------------------------|---------------|
| TEMPERATURE: Thermometer ID: SC2 (Criteria: 0.0 °C – 6.0 °C, not froze | n except se | diment/tissu | e) |
| Temperature 3.5°C - 0.2°C (CF) = 3.3°C | Blank | ☐ Sample | е |
| ☐ Sample(s) outside temperature criteria (PM/APM contacted by:) | | | |
| ☐ Sample(s) outside temperature criteria but received on ice/chilled on same c | lay of sampl | ing. | |
| ☐ Received at ambient temperature, placed on ice for transport by Co | | | |
| Ambient Temperature: ☐ Air ☐ Filter | | Checked b | v: 15 |
| | | | |
| CUSTODY SEALS INTACT: | | | 1== |
| ☑Cooler □ □ No (Not Intact) □ Not Present | □ N/A | | |
| ☑Sample □ □ No (Not Intact) □ Not Present | | Checked by | y: <u>836</u> |
| CANDIT CONDITION. | V | NO. | NI/Λ |
| | Yes | No | N/A |
| Chain-Of-Custody (COC) document(s) received with samples | | | |
| COC document(s) received complete | | | |
| ☐ Collection date/time, matrix, and/or # of containers logged in based on sample labels. | | | |
| □ No analysis requested. □ Not relinquished. □ No date/time relinquished. | | П | |
| Sampler's name indicated on COC | | | |
| Sample container label(s) consistent with COC | | <u></u> | |
| Sample container(s) intact and good condition | • | <u></u> | |
| Proper containers and sufficient volume for analyses requested | | | U |
| Analyses received within holding time | | in northead | Ц |
| Aqueous samples received within 15-minute holding time | | | |
| ☐ pH ☐ Residual Chlorine ☐ Dissolved Sulfides ☐ Dissolved Oxygen | | | |
| Proper preservation noted on COC or sample container | | | |
| ✓ Unpreserved vials received for Volatiles analysis | ر | _ | |
| Volatile analysis container(s) free of headspace | | | |
| Tedlar bag(s) free of condensation CONTAINER TYPE: | . 🗆 | | |
| Solid: □4ozCGJ □8ozCGJ □16ozCGJ □Sleeve () □EnCore | s® □Terra | Cores [®] □_ | |
| Aqueous: ZVOA ZVOAh □VOAna₂ □125AGB □125AGBh □125AGBp | | | |
| □500AGB Ø500AGJ □500AGJs □250AGB □250CGB □250CGBs | 21PBn | □1PB na □ | 1500PB |
| □250PB □250PBn □125PB □125PB znna □100PJ □100PJ na₂ □ | | | |
| Air: □Tedlar® □Canister Other: □ Trip Blank Lot#: 140 90 81 | | | |
| Container: C: Clear A: Amber P: Plastic G: Glass J: Jar B: Bottle Z: Ziploc/Resealable Bag E: Er Preservative: h: HCL n: HNO ₃ na ₂ :Na ₂ S ₂ O ₃ na: NaOH p: H ₃ PO ₄ s: H ₂ SO ₄ u: Ultra-pure znna: ZnAc ₂ +Na | • | Reviewed by: Scanned by | |



Supplemental Report 1

Calscience



WORK ORDER NUMBER: 14-10-2234

The difference is service



AIR SOIL WATER | MARINE CHEMISTRY

Analytical Report For

Client: Environmental Science International, Inc.

Client Project Name: Red Hill LTM 112066

Attention: Jeff Hattemer

354 Uluniu Street, Suite 304 Kailua, HI 96734-2500

Rand Ollo

Approved for release on 11/05/2014 by: Richard Villafania Project Manager



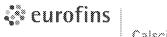
Resultance

Emailyour PM &

Eurofins Calscience, Inc. (Calscience) certifies that the test results provided in this report meet all NELAC requirements for parameters for which accreditation is required or available. Any exceptions to NELAC requirements are noted in the case narrative. The original report of subcontracted analyses, if any, is attached to this report. The results in this report are limited to the sample(s) tested and any reproduction thereof must be made in its entirety. The client or recipient of this report is specifically prohibited from making material changes to said report and, to the extent that such changes are made, Calscience is not responsible, legally or otherwise. The client or recipient agrees to indemnify Calscience for any defense to any litigation which may arise.

7440 (mergis Vice Contours) core (A.9264) (437 × TEL-710 895-599) - FAX (714 896-790) - programmer con-

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Calscience

Contents

| Client Pro | ject Name: | Red | IIIH b | LTM 11 | 12066 | | | | | | | | | | | | | | | | |
|------------|--|-------------|----------|--------|-------|-------|-------|-----|------|----|------|----------|------|------|------|----------|------|------|----------|--|------------------|
| Work Ord | er Number: | 14- | 10-22 | 34 | | | | | | | | | | | | | | | | | |
| 1 | Work Ord | ler | Nar | rativ | e | | | | | | | | | | | | | | | | 3 |
| 2 | Client Sar 2.1 EPA | | | | | | | | | | | | | | | | | | | | 4 4 |
| 3 | Quality Co 3.1 MS/N 3.2 PDS/ 3.3 LCS/ | /ISI /PC | D DSD | | | | | | | | | | | | | | | | | | 5 5 6 7 |
| 4 | Sample A | ۱na | lysi | s Sui | mma | ıry. | | | | | | | | | | | | | | | 8 |
| 5 | Glossary | of | Teri | ns a | nd C |)uali | ifier | S | | | | | | | | | | | | | 9 |
| 6 | Chain-of- | Cu | stoc | ly/Sa | ampl | e Re | ece | ipt | Foi | m. | | | | | | | | | | | 10 |



Work Order Narrative

Work Order: 14-10-2234 Page 1 of 1

Condition Upon Receipt:

Samples were received under Chain-of-Custody (COC) on 10/29/14. They were assigned to Work Order 14-10-2234.

Unless otherwise noted on the Sample Receiving forms all samples were received in good condition and within the recommended EPA temperature criteria for the methods noted on the COC. The COC and Sample Receiving Documents are integral elements of the analytical report and are presented at the back of the report.

Holding Times:

All samples were analyzed within prescribed holding times (HT) and/or in accordance with the Calscience Sample Acceptance Policy unless otherwise noted in the analytical report and/or comprehensive case narrative, if required.

Any parameter identified in 40CFR Part 136.3 Table II that is designated as "analyze immediately" with a holding time of <= 15 minutes (40CFR-136.3 Table II, footnote 4), is considered a "field" test and the reported results will be qualified as being received outside of the stated holding time unless received at the laboratory within 15 minutes of the collection time.

Quality Control:

All quality control parameters (QC) were within established control limits except where noted in the QC summary forms or described further within this report.

Additional Comments:

Air - Sorbent-extracted air methods (EPA TO-4A, EPA TO-10, EPA TO-13A, EPA TO-17): Analytical results are converted from mass/sample basis to mass/volume basis using client-supplied air volumes.

New York NELAP air certification does not certify for all reported methods and analytes, reference the accredited items here: http://www.calscience.com/PDF/New_York.pdf

Solid - Unless otherwise indicated, solid sample data is reported on a wet weight basis, not corrected for % moisture. All QC results are always reported on a wet weight basis.

Subcontractor Information:

Unless otherwise noted below (or on the subcontract form), no samples were subcontracted.



Analytical Report

Environmental Science International, Inc.

Date Received:

Work Order:

10/29/14

Work Order:

14-10-2234

Kailua, HI 96734-2500

Preparation:

Method:

Units:

Units:

Units:

10/29/14

Project: Red Hill LTM 112066 Page 1 of 1

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|----------------------|----------------------|------------------------|--------|------------|------------------|-----------------------|-------------|
| ES117 | 14-10-2234-2-J | 10/28/14 12:00 | | ICP/MS 04 | 10/31/14 | 11/03/14 14:46 | 141031L01D |

Comment(s): - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag.

Parameter

Lead - Result - R

| Method Blank | 099-16-094- | 581 N/A | Aqueous IC | P/MS 04 10/31/ | 14 11/03/14 14:31 | 141031L01D |
|------------------|---|-----------------------|--------------------|------------------------|------------------------|-------------|
| Comment(s): | - Results were evaluated to the MDL (DL | .), concentrations >= | to the MDL (DL) bu | it < RL (LOQ), if four | nd, are qualified with | a "J" flag. |
| <u>Parameter</u> | | Result | <u>RL</u> | MDL. | <u>DF</u> | Qualifiers |
| Lead | | <0.0898 | 1.00 | 0.0898 | 1.00 | U |

RL: Reporting Limit. DF: Dilution Factor. MDL: Method Detection Limit.



Quality Control - Spike/Spike Duplicate

Environmental Science International, Inc.

Date Received:

Work Order:

14-10-2234

Kailua, HI 96734-2500

Preparation:

N/A

Method:

EPA 200.8

Project: Red Hill LTM 112066 Page 1 of 1

| Quality Control Sample ID | Туре | | Matrix | In | strument | Date Prepared | Date Ana | lyzed | MS/MSD Bat | tch Number |
|---------------------------|-----------------|-----------------------|-------------|--------------------|--------------|---------------|----------|-------|------------|-------------------|
| ES117 | Sample | | Aqueous | . IC | P/MS 04 | 10/31/14 | 11/03/14 | 14:46 | 141031S01A | |
| ES117 | Matrix Spike | | Aqueous | i ic | P/MS 04 | 10/31/14 | 11/03/14 | 14:35 | 141031S01A | |
| ES117 | Matrix Spike Du | plicate | Aqueous | i ic | P/MS 04 | 10/31/14 | 11/03/14 | 14:37 | 141031S01A | i. |
| Parameter | | <u>Spike</u> Added | MS Conc. | <u>MS</u> %Rec. | MSD Conc. | MSD %Rec. | %Rec. CL | RPD | RPD CL | <u>Qualifiers</u> |
| Lead | ND 1 | 100.0 | 100.7 | 101 | 100.8 | 101 | 80-120 | 0 | 0-20 | |





Quality Control - PDS

Environmental Science International, Inc.

Date Received:

Work Order:

10/29/14

Work Order:

Kailua, HI 96734-2500

Preparation:

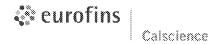
N/A

Method: EPA 200.8

Project: Red Hill LTM 112066 Page 1 of 1

| Quality Control Sample ID | Type | M | //atrix | Instrument | Date Prepared [| | DS/PDSD Batch umber | |
|---------------------------|--------|--------------|-------------|------------|-----------------|------------------|------------------------|--|
| ES117 | Sample | Д | Aqueous | ICP/MS 04 | 10/31/14 00:00 | 11/03/14 14:46 1 | 41031S01A | |
| ES117 | PDS | Д | vqueous | ICP/MS 04 | 10/31/14 00:00 | 11/03/14 14:39 1 | 41031S01A | |
| Parameter | | Sample Conc. | Spike Added | PDS Conc | . PDS %Rec | . %Rec. CL | <u>Qualifiers</u> | |
| Lead | | ND | 100.0 | 99.82 | 100 | 75-125 | | |





Quality Control - LCS

Environmental Science International, Inc.

Date Received:

Work Order:

14-10-2234

Kailua, HI 96734-2500

Preparation:

N/A

Method:

EPA 200.8

Project: Red Hill LTM 112066 Page 1 of 1

| Quality Control Sample ID Type | | Matrix | Instrument Da | , | , | LCS Batch Number |
|--------------------------------|-----|-------------|-----------------|-------------|------------|------------------|
| 099-16-094-581 | LCS | Aqueous | ICP/MS 04 10 | /31/14 11/0 | 3/14 14:33 | 141031L01D |
| <u>Parameter</u> | | Spike Added | Conc. Recovered | LCS %Rec. | %Rec | . CL Qualifiers |
| Lead | | 100.0 | 95.21 | 95 | 80-120 |) |

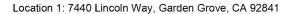






Sample Analysis Summary Report

| Work Order: 14-10-2234 | | | | Page 1 of 1 |
|------------------------|------------|------------|-------------------|---------------------|
| Method | Extraction | Chemist ID | <u>Instrument</u> | Analytical Location |
| EPA 200.8 | N/A | 598 | ICP/MS 04 | 1 |





Glossary of Terms and Qualifiers

Work Order: 14-10-2234 Page 1 of 1

| *************************************** | 5 |
|---|--|
| Qualifiers | <u>Definition</u> |
| * | See applicable analysis comment. |
| < | Less than the indicated value. |
| > | Greater than the indicated value. |
| 1 | Surrogate compound recovery was out of control due to a required sample dilution. Therefore, the sample data was reported without further clarification. |
| 2 | Surrogate compound recovery was out of control due to matrix interference. The associated method blank surrogate spike compound was in control and, therefore, the sample data was reported without further clarification. |
| 3 | Recovery of the Matrix Spike (MS) or Matrix Spike Duplicate (MSD) compound was out of control due to suspected matrix interference. The associated LCS recovery was in control. |
| 4 | The MS/MSD RPD was out of control due to suspected matrix interference. |
| 5 | The PDS/PDSD or PES/PESD associated with this batch of samples was out of control due to suspected matrix interference. |
| 6 | Surrogate recovery below the acceptance limit. |
| 7 | Surrogate recovery above the acceptance limit. |
| В | Analyte was present in the associated method blank. |
| BU | Sample analyzed after holding time expired. |
| BV | Sample received after holding time expired. |
| E | Concentration exceeds the calibration range. |
| ET | Sample was extracted past end of recommended max. holding time. |
| HD | The chromatographic pattern was inconsistent with the profile of the reference fuel standard. |
| HDH | The sample chromatographic pattern for TPH matches the chromatographic pattern of the specified standard but heavier hydrocarbons were also present (or detected). |
| HDL | The sample chromatographic pattern for TPH matches the chromatographic pattern of the specified standard but lighter hydrocarbons were also present (or detected). |
| J | Analyte was detected at a concentration below the reporting limit and above the laboratory method detection limit. Reported value is estimated. |
| JA | Analyte positively identified but quantitation is an estimate. |
| ME | LCS Recovery Percentage is within Marginal Exceedance (ME) Control Limit range (+/- 4 SD from the mean). |
| ND | Parameter not detected at the indicated reporting limit. |
| Q | Spike recovery and RPD control limits do not apply resulting from the parameter concentration in the sample exceeding the spike concentration by a factor of four or greater. |
| | |

- SG The sample extract was subjected to Silica Gel treatment prior to analysis.
- U Undetected at the laboratory method detection limit.
- X % Recovery and/or RPD out-of-range.
- Z Analyte presence was not confirmed by second column or GC/MS analysis.

Solid - Unless otherwise indicated, solid sample data is reported on a wet weight basis, not corrected for % moisture. All QC results are reported on a wet weight basis.

Any parameter identified in 40CFR Part 136.3 Table II that is designated as "analyze immediately" with a holding time of <= 15 minutes (40CFR-136.3 Table II, footnote 4), is considered a "field" test and the reported results will be qualified as being received outside of the stated holding time unless received at the laboratory within 15 minutes of the collection time.

A calculated total result (Example: Total Pesticides) is the summation of each component concentration and/or, if "J" flags are reported, estimated concentration. Component concentrations showing not detected (ND) are summed into the calculated total result as zero concentrations.

| , | | | | | | | | | | | | | www.compressure. | | | Page | 10 c | of 12 |
|--|-------------------------------|-------------------------------------|--------------|---|--|--|--|-----------------------------|---------------------|--|--------------------|------------|--|---------------------------|-----------------------------|---|--|--------------------------------------|
| | | UNT) | Ę | | | | (8'001 | ·*/ 3 | רפטי רפטי | NAT TO THE PROPERTY OF THE PRO | × | × | | | | Time: \S\chi_VS\chiVS\chi_VS\chi_VS\chi_VS\chi_VS\chi_VS\chi_VS\chi_VS\chi_VS\c | Time: | Time: |
| 78/16 1 or 1 | P.O. NO.: | SAMPLER(S); (PRINT) | 以市局 | ANALYSES | as needed. | | MIS 0728 18 □ X747\011 18.812 □ 9917 | X 2 0728 09 □ el£ | TS2 Meta | Holodonic control de contraction de | X | X | | | | Date: | Ďaře: | Date; |
| Date C | R: | 35 | Ž. | REQUESTED A | check box or fill in blank as n | eroD s | ore 🗆 Tema | | Prep (50 | | | | | | | | | Salkt |
| 0#/LAB USE ONLY | CLIENT PROJECT NAME / NUMBER: | ROJECT CONTACT: | Hattomer | Y | Please check box | | 85e0 ==================================== | □ 38T <i>N</i> | HqT | × | 8 | X | | | | ansoning and the contract of t | | |
| | CLIENT PRC | PROJECT O | C 8 | _ <u>こ</u> | | * | | tered (1) | | × | X | × | The second secon | | | Received by: (Signature/Affiliation) | Received by: (Signature/Affiliation) | Received by: (Signature/Affiliation) |
| 5-5494 us Christi | | | 734 ZIP. | F-MAIL: Shattemer Clescience, com, Albert Perience, com | JARD | LOG CODE | | | Unpreserv | × | W. | У | 100 000 000 000 000 000 000 000 000 000 | near agratus and agration | noodaa vaassankoodaa qahasa | Received by: (Si | Received by: (Si | Received by: (Si |
| -1427 • (714) 89 ouston, and Corp off information, | | | STATE: 967 | Com Algre | RD"): .YS (XSTANDARD | Andrew missing missing management and an analysis of the property of the prope | Holist | | MATRIX OF CONT. | No. | \$ \$ \(\times \) | i i | | | | X | | |
| Nay, Garden Grove, CA 92841-1427 • (7 Concord, San Luis Obispo, Houston, and courier service / sample drop off informati contact sales@calscience.com or call us. | | 1 to 304 | j | COSCIENCE | o any TAT not "STANDARD" 72 HR | | ring anal | | SAMPLING TE TIME | 0011 41/82/01 | 0001 1000 | S/W 114750 | | | | | | |
| 7440 Lincoln Way, Garden Grove, CA 92841-1427。 (714) 895-5494 Other locations: Concord, San Luis Obispo, Houston, and Corpus Christi For courier service / sample drop off information, contact sales@calscience.com or call us. | | スプスマ | | E-MAIL: | ges may apply to any □ 48 HR □ 7 | emochemistics and amochemistics from the control of | of month | | S DATE | 202 | | 10/26/14 | | | | | The state of the s | |
| 7440 Lino. | LABORATORY CLIENT: 651 | ADDRESS: 354 Ulyning St., Suite 304 | OITY: Kailya | 7EL (808) 261-0740 | TURNAROUND TIME (Rush surcharges may apply to any TAT not "STANDARD"): SAME DAY 124 HR 5 DAYS | COELT EDF GLOBAL ID | special instructions: Degular (full) fim monitaring analytelist | | SAMPLEID | 65TRIP | から | MS-18 | | | | Relinquished by: (Signature) | Relinquished by: (Signature) | Relinquished by: (Signature) |
| | BOF | ODRE | ТҮ: | | JRN/ JSAN | 100 |) PEC. | | LAB USE | 1 | N | 2 | | | | | eill G | elinc |

ORIGIN ID:HNLA

SHIP DATE: 280CT14 ACTWGT: 52.5 LB CAD: /POS1525 DIMS: 26x15x14 IN

BILL RECIPIENT

UNITED STATES US

TO

CALSCIENCE 7440 LINCOLN WAY

GARDEN GROVE CA 92841

(714) 895-5494

REF:

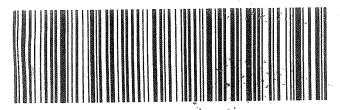


TRK# 8045 5791 7375

WED - 29 OCT AA STANDARD OVERNIGHT

WZ APVA

92841 ca=us SNA



(223⁴)



Calscience

WORK ORDER #: 14-10-223 [4

SAMPLE RECEIPT FORM

M Cooler <u>/</u> of <u>/</u>

| CLIENT: <u>ESI</u> DATE: _ | 10/27 | / 17 |
|--|----------------------|--|
| TEMPERATURE: Thermometer ID: SC2 (Criteria: 0.0 °C – 6.0 °C, not frozen except sec | diment/tiss | ue) |
| Temperature 3 .5 °C - 0.2 °C (CF) = 3 .3 °C ☑ Blank | ☐ Samp | le |
| ☐ Sample(s) outside temperature criteria (PM/APM contacted by:) | · | |
| ☐ Sample(s) outside temperature criteria but received on ice/chilled on same day of sampli | na. | |
| ☐ Received at ambient temperature, placed on ice for transport by Courier. | | |
| Ambient Temperature: Air Filter | Checked I | hv: 15 |
| Ambient reinperature. El All El Filtor | | ~ J · |
| CUSTODY SEALS INTACT: | | £ - |
| ☑Cooler □ □ No (Not Intact) □ Not Present □ N/A | Checked b | ру: |
| ☑Sample □ □ No (Not Intact) □ Not Present | Checked b | ру: <u>836</u> |
| | | |
| SAMPLE CONDITION: Yes | No — | N/A |
| Chain-Of-Custody (COC) document(s) received with samples | | |
| COC document(s) received complete | | |
| ☐ Collection date/time, matrix, and/or # of containers logged in based on sample labels. | | |
| ☐ No analysis requested. ☐ Not relinquished. ☐ No date/time relinquished. | | |
| Sampler's name indicated on COC | | |
| Sample container label(s) consistent with COC | | |
| Sample container(s) intact and good condition | | |
| Proper containers and sufficient volume for analyses requested | | |
| Analyses received within holding time | | |
| Aqueous samples received within 15-minute holding time | | |
| □ pH □ Residual Chlorine □ Dissolved Sulfides □ Dissolved Oxygen □ | | |
| Proper preservation noted on COC or sample container | | |
| Unpreserved vials received for Volatiles analysis | | |
| Volatile analysis container(s) free of headspace | | |
| Tedlar bag(s) free of condensation CONTAINER TYPE: | | 4 |
| Solid: □4ozCGJ □8ozCGJ □16ozCGJ □Sleeve () □EnCores® □Terra | Cores [®] 🗆 | MA MANANA MA |
| Aqueous: ØVOA ØVOAh □VOAna₂ □125AGB □125AGBh □125AGBp Ø1AGB □ |]1AGB na ₂ | □1AGB s |
| □500AGB Ø500AGJ □500AGJs □250AGB □250CGB □250CGBs Ø1PBn | | |
| □250PB □250PBn □125PB □125PB znna □100PJ □100PJ na₂ □ □ | | |
| Air: Tedlar® Canister Other: Trip Blank Lot#: 140908B Labeled/Container: C: Clear A: Amber P: Plastic G: Glass J: Jar B: Bottle Z: Ziploc/Resealable Bag E: Envelope R Preservative: h: HCL n: HNO ₃ na ₂ :Na ₂ S ₂ O ₃ na: NaOH p: H ₃ PO ₄ s: H ₂ SO ₄ u: Ultra-pure znna: ZnAc ₂ +NaOH f: Filtered | eviewed by | : <u>852</u> |

RAW DATA SHEET FOR METHOD: EPA 8015B (M)

WORK ORDER: **INSTRUMENT:**

EXTRACTION:

14-10-2234

GC 45

EPA 3510C

D/T EXTRACTED: 2014-10-30 00:00

ANALYZED BY:

D/T ANALYZED:

2014-10-31 11:16

682

REVIEWED BY: D/T REVIEWED:

DATA FILE:

W:\GC 45\GC 45 DATA\2014\141030\14103067.D\14103067

2 #

CLIENT SAMPLE NUMBER: ES117

LCS/MB BATCH:

141030B06

SAMPLE VOLUME / WEIGHT:

DEFAULT: 500.00 ml / ACTUAL: 500.00 ml

MS/MSD BATCH: FINAL VOLUME / WEIGHT: DEFAULT: 5.00 ml / ACTUAL: 2.50 ml

LOD

12

ADJUSTMENT RATIO TO PF: UNITS: ug/L

0.50

COMMENT:

Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, as

qualified with a "J" flag. TPH as Diesel is quantified in the carbon range C10-C28.

COMPOUND

INI. CONC DF

1.00

CONC

22.2

DL

LOQ

QUAL

TPH as Diesel

4440

11

25

bJ

Area Percent Report

Data File Name : W:\GC_45\GC 45 DATA\2014\141030\14103067.D

Vial Number : Vial 67

Page Number : 1
Operator : 682
Instrument : GC 45 Injection Number: 1 Sequence Line : 67 Sample Name : 14-10-2234-2

Instrument Method: C:\CHEM32\1\METHODS\ ->

Acquired on : 31 Oct 14 11:16 am

Report Created on: 04 Nov 14 02:51 pm Analysis Method : 8015B.MTH

Software Revision: Rev. B.03.02 [341] Copyright © Agilent Technologies

Sig. 1 in W:\GC 45\GC 45 DATA\2014\141030\ ->

| Pk | Ret Time | Area | - | | | Response % |
|----|----------|--------|-----|----|-------|------------|
| 1 | | | | | | 1 |
| 1 | 2.455 | 0.77 | 0 | VV | 0.054 | 0.105 |
| 2 | 2.637 | 1.41 | 0 | VV | 0.039 | 0.193 |
| 3 | 2.694 | 1.80 | 1 | VV | 0.051 | 0.247 |
| 4 | 2.750 | 2.10 | 1 | VV | 0.036 | 0.287 |
| 5 | 2.837 | 0.86 | 0 | VV | 0.032 | 0.118 |
| 6 | 2.885 | 0.68 | 0 | VV | 0.024 | 0.093 |
| 7 | 2.945 | 1.24 | 0 | VV | 0.052 | 0.170 |
| 8 | 3.169 | 5.53 | 1 | VV | 0.053 | 0.756 |
| 9 | 3.272 | 2.00 | 1 | VV | 0.044 | 0.273 |
| 10 | 3.344 | 4.42 | 1 | VV | 0.067 | 0.605 |
| 11 | 3.437 | 3.47 | 1 | VV | 0.044 | 0.474 |
| 12 | 3.486 | 11.09 | 1 | VV | 0.130 | 1.517 |
| 13 | 3.718 | 7.85 | 1 | VV | 0.107 | 1.074 |
| 14 | 3.850 | 2.34 | 1 | VV | 0.041 | 0.321 |
| 15 | 3.916 | 2.44 | 1 | VV | 0.054 | 0.333 |
| 16 | 4.020 | 2.87 | 2 | VV | 0.024 | 0.393 |
| 17 | 4.082 | 2.19 | 0 | VB | 0.081 | 0.300 |
| 18 | 4.239 | 2.57 | 1 | BV | 0.056 | 0.351 |
| 19 | 4.403 | 1.51 | 0 | VV | 0.049 | 0.206 |
| 20 | 4.444 | 1.36 | 0 | VB | 0.042 | 0.186 |
| 21 | 4.590 | 0.91 | 0 | BV | 0.061 | 0.125 |
| 22 | 4.701 | 1.77 | 0 | VV | 0.067 | 0.242 |
| 23 | 4.850 | 0.97 | 0 | VV | 0.033 | 0.132 |
| 24 | 4.911 | 0.53 | 0 | VV | 0.030 | 0.072 |
| 25 | 4.967 | 0.44 | 0 | VV | 0.028 | 0.061 |
| 26 | 5.069 | 1.34 | 1 | VV | 0.027 | 0.184 |
| 27 | 5.376 | 666.87 | 439 | VV | 0.025 | 91.184 |

Total area = 731.35

Area Percent Report

Data File Name : W:\GC_45\GC 45 DATA\2014\141030\14103067.D

Page Number : 2

Operator : 682 Vial Number : Vial 67

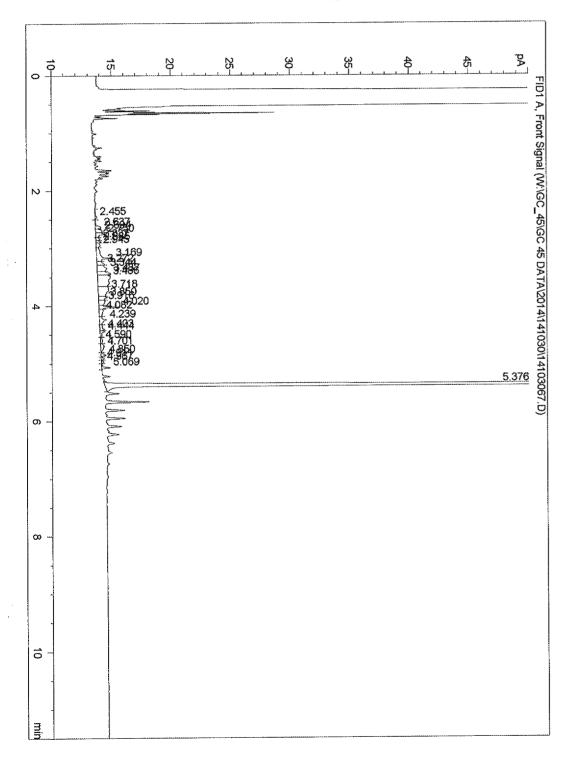
Instrument : GC 45 Injection Number : 1
Sample Name : 14-10-2234-2 Sequence Line : 67

Instrument Method: C:\CHEM32\1\METHODS\ ->

Acquired on : 31 Oct 14 11:16 am

Report Created on: 04 Nov 14 02:51 pm Analysis Method : 8015B.MTH

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RAW DATA SHEET FOR METHOD: EPA 8015B (M)

WORK ORDER: INSTRUMENT:

14-10-2234

GC 45

EXTRACTION: D/T EXTRACTED: 2014-10-30 00:00

DATA FILE:

EPA 3510C

ANALYZED BY:

D/T ANALYZED:

682

2014-10-31 11:35

REVIEWED BY: D/T REVIEWED:

W:\GC 45\GC 45 DATA\2014\141030\14103068.D\14103068

3

CLIENT SAMPLE NUMBER: ES118

LCS/MB BATCH: MS/MSD BATCH:

141030B06

SAMPLE VOLUME / WEIGHT:

DEFAULT: 500.00 ml / ACTUAL: 500.00 ml

DEFAULT: 5.00 ml / ACTUAL: 2.50 ml FINAL VOLUME / WEIGHT:

UNITS:

ug/L

ADJUSTMENT RATIO TO PF:

0.50

11

COMMENT:

Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, as

qualified with a "J" flag. TPH as Diesel is quantified in the carbon range C10-C28.

COMPOUND TPH as Diesel INI. CONC DF

3240

1.00

CONC

16.2

LOD <u>DL</u> 12

<u>LOQ</u> 25

QUAL bJ

Area Percent Report

Data File Name : W:\GC_45\GC 45 DATA\2014\141030\14103068.D

Page Number : 1

Operator : 682 Vial Number : Vial 68

Instrument : GC 45 Injection Number : 1 Sample Name : 14-10-2234-3 Sequence Line : 68

Instrument Method: C:\CHEM32\1\METHODS\ ->

Acquired on : 31 Oct 14 11:35 am

Report Created on: 04 Nov 14 02:53 pm Analysis Method : 8015B.MTH

Software Revision: Rev. B.03.02 [341] Copyright © Agilent Technologies

Sig. 1 in W:\GC 45\GC 45 DATA\2014\141030\ ->

| Pk | Ret Time | Area | • | | Width | Response % |
|--------|----------|----------------------------------|-----|----|-------|------------|
| • | | +++ +++ +++ ++ + +- +- +- | | | | |
| 1 | 2.457 | 0.55 | | VV | 0.050 | 0.091 |
| 2 | 2.601 | 0.48 | | VV | 0.021 | 0.080 |
| 3 | 2.640 | 0.60 | | VV | 0.024 | 0.101 |
| 4 | 2.689 | 1.65 | | VV | 0.058 | 0.276 |
| 5 | 2.752 | 1.54 | | VV | 0.038 | 0.257 |
| 6 | 2.837 | 0.65 | | VV | 0.034 | 0.108 |
| 7 | 2.885 | 0.64 | 0 | VV | 0.033 | 0.107 |
| 8 | 2.942 | 0.65 | 0 | VV | 0.043 | 0.109 |
| 9 | 3.101 | 1.29 | 0 | VV | 0.073 | 0.216 |
| 10 | 3.169 | 3.26 | 2 | VV | 0.026 | 0.545 |
| 11 | 3.342 | 3.49 | 1 | VV | 0.110 | 0.583 |
| 12 | 3.437 | 1.96 | 1 | VV | 0.049 | 0.327 |
| 13 | 3,481 | 1.65 | 1 | VV | 0.040 | 0.275 |
| 14 | 3.527 | 1.70 | 1 | VV | 0.042 | 0.284 |
| 15 | 3.596 | 3.74 | 1 | VV | 0.095 | 0.624 |
| 16 | 3.684 | 1.55 | 1 | VV | 0.047 | 0.259 |
| 17 | 3.755 | 3.47 | 1 | VV | 0.089 | 0.579 |
| 18 | 3.851 | 1.60 | 1 | VV | 0.045 | 0.268 |
| 19 | 3.916 | 0.61 | 0 | VV | 0.024 | 0.103 |
| 20 | 3.944 | 1.04 | 0 | VV | 0.047 | 0.174 |
| 21 | 4.020 | 2.38 | 2 | VV | 0.022 | 0.398 |
| 22 | 4.089 | 1.00 | 0 | VV | 0.054 | 0.167 |
| 23 | 4.132 | 0.32 | 0 | VV | 0.024 | 0.053 |
| 24 | 4.160 | 0.20 | 0 | VV | 0.022 | 0.033 |
| 25 | 4.239 | 2.07 | 1 | VV | 0.056 | 0.346 |
| 26 | 4.357 | 0.79 | 0 | VV | 0.044 | 0.132 |
| 27 | 4.406 | 1.51 | 0 | VV | 0.055 | 0.253 |
| 28 | 4.596 | 0.74 | 0 | VV | 0.065 | 0.124 |
| 29 | 4.723 | 1.18 | 0 | VV | 0.054 | 0.198 |
| 30 | 4.737 | 1.47 | 1 | VV | 0.030 | 0.246 |
| 31 | 4.851 | 0.96 | 0 | VV | 0.037 | 0.161 |
| 32 | 4.913 | 0.71 | 0 | VV | 0.030 | 0.118 |
| 33 | 4.972 | 0.41 | 0 | VV | 0.023 | 0.069 |
| 34 | 5.024 | 0.16 | 0 | VV | 0.023 | 0.027 |
| 35 | 5.073 | 0.91 | 1 | VV | 0.024 | 0.153 |
| 36 | 5.270 | 0.06 | 0 | VM | 0.010 | 0.010 |
| 37 | 5,379 | 551.80 | 370 | VV | 0.025 | 92.147 |

Area Percent Report

Data File Name : W:\GC_45\GC 45 DATA\2014\141030\14103068.D

Page Number : 2

Vial Number : Vial 68 : 682

Operator : 682
Instrument : GC 45 Injection Number: 1 Sample Name : 14-10-2234-3 Sequence Line : 68

Instrument Method: C:\CHEM32\1\METHODS\ ->

Acquired on : 31 Oct 14 11:35 am

Report Created on: 04 Nov 14 02:53 pm Analysis Method : 8015B.MTH

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Total area = 598.83

Area Percent Report

Data File Name : W:\GC_45\GC 45 DATA\2014\141030\14103068.D

Page Number : 3

Operator : 682 Vial Number : Vial 68

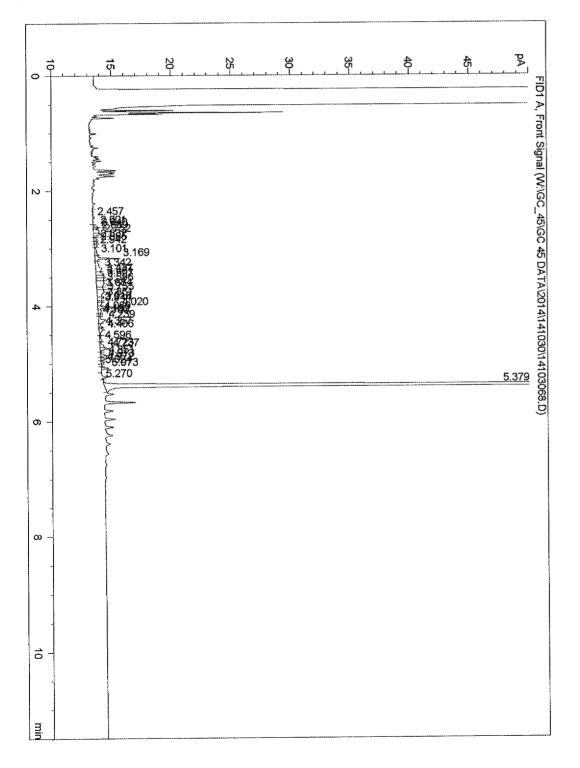
Instrument : GC 45 Injection Number : 1
Sample Name : 14-10-2234-3 Sequence Line : 68

Instrument Method: C:\CHEM32\1\METHODS\ ->

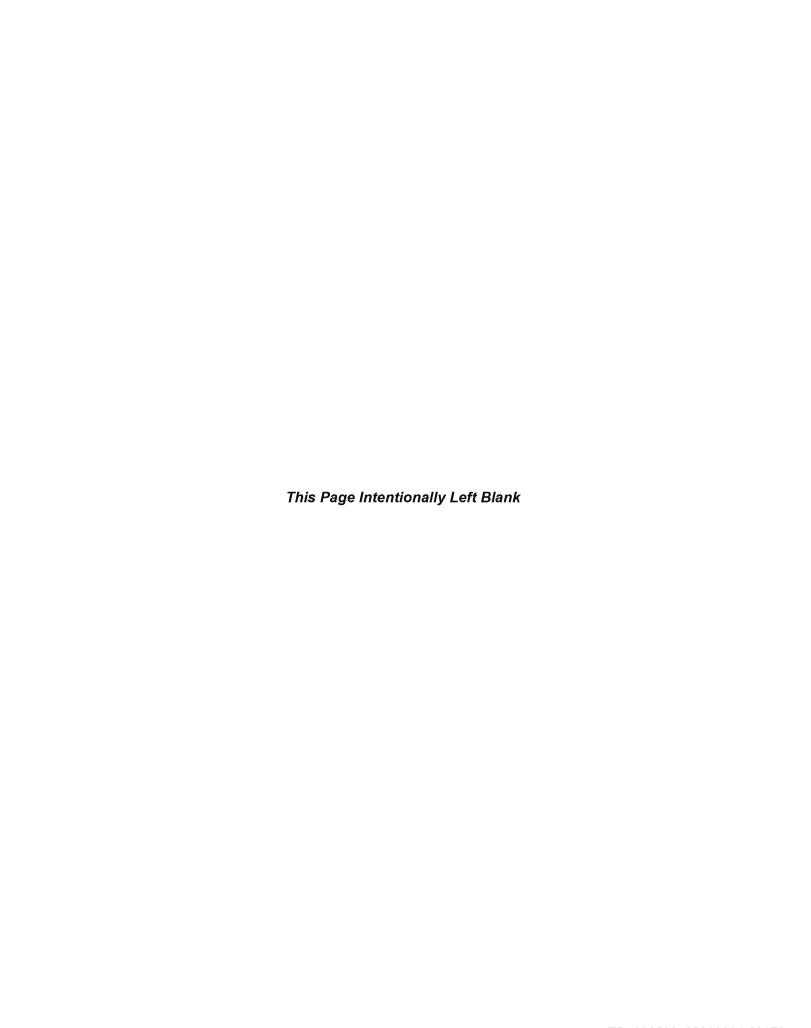
Acquired on : 31 Oct 14 11:35 am

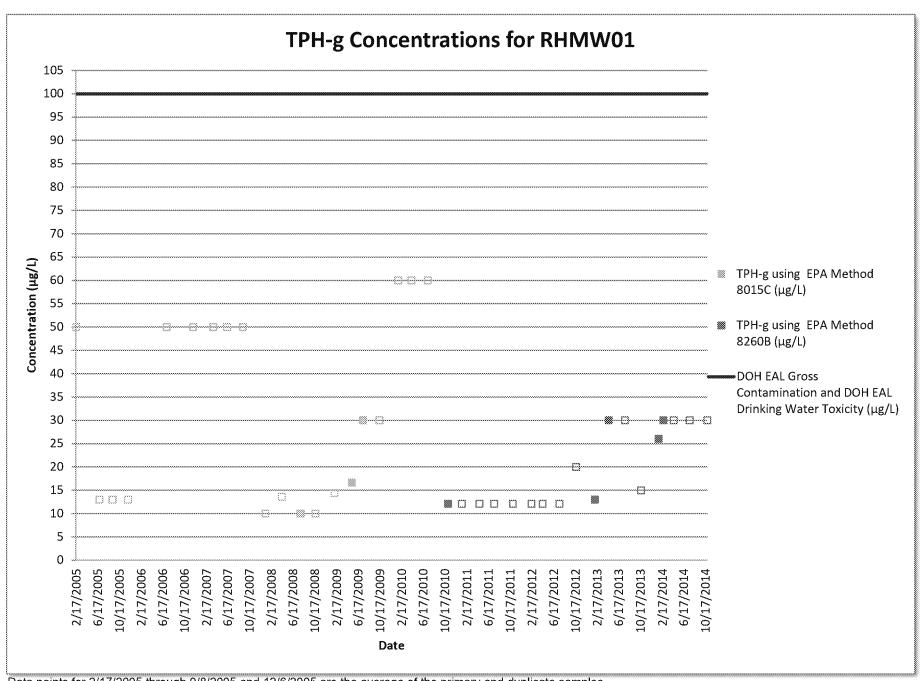
Report Created on: 04 Nov 14 02:53 pm Analysis Method : 8015B.MTH

Software Revision: Rev. B.03.02 [341] Copyright © Agilent Technologies



APPENDIX D Historical Groundwater Exceedances

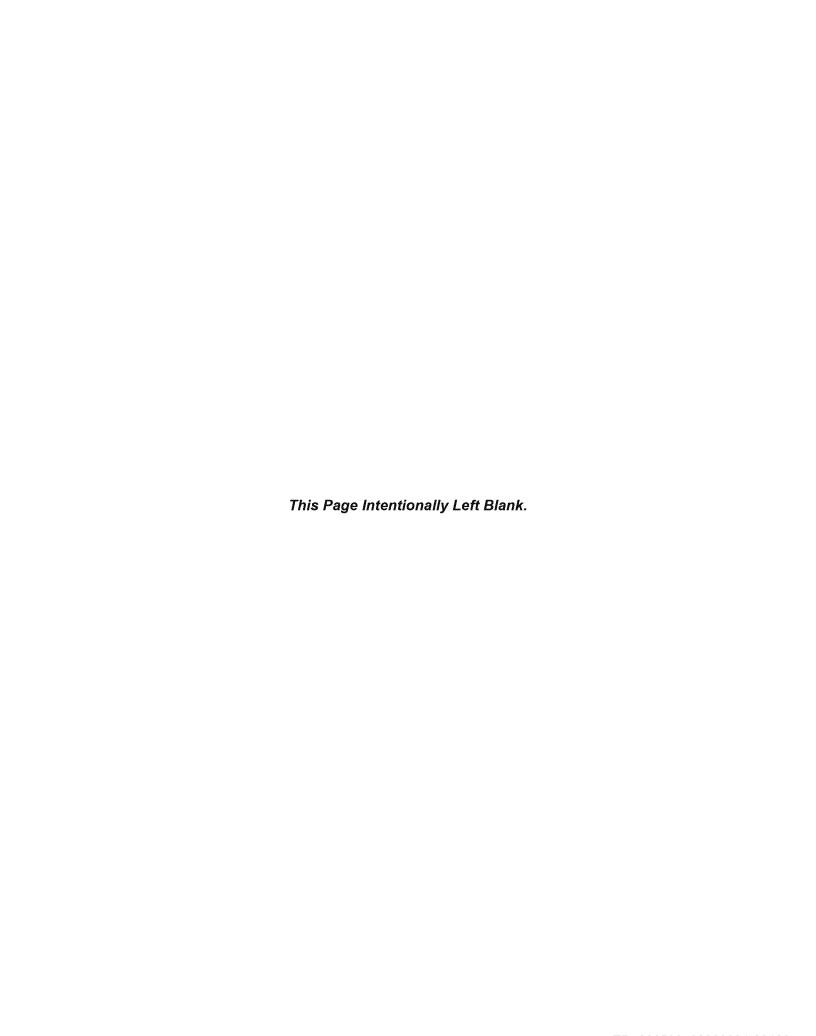


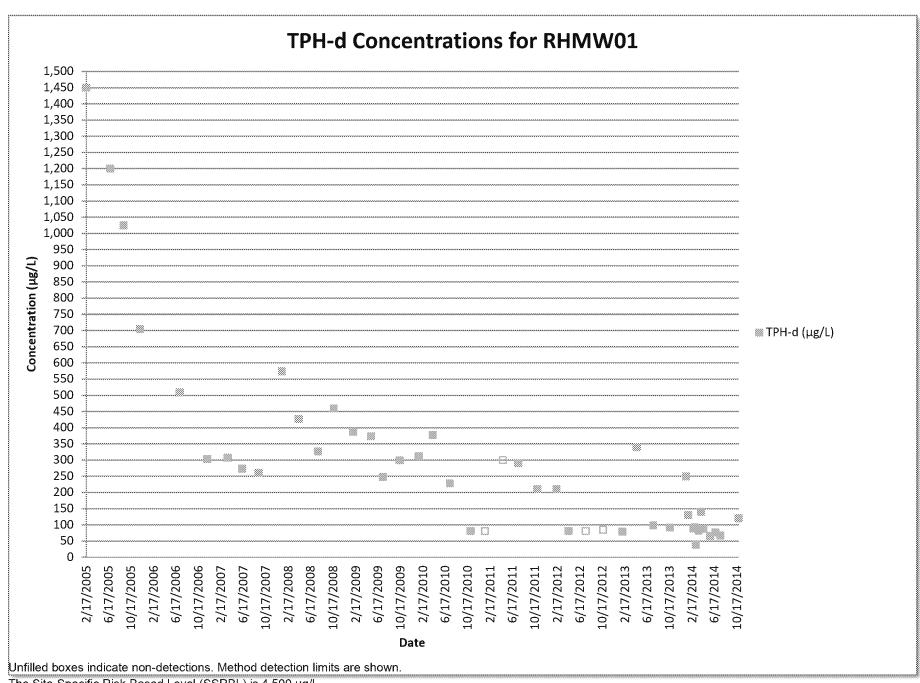


Data points for 2/17/2005 through 9/8/2005 and 12/6/2005 are the average of the primary and duplicate samples.

Possible laboratory contamination for 10/22/2012, 10/21/2013, and 1/28/2014 sampling events.

Unfilled boxes indicate non-detections. Method detection limits are shown.

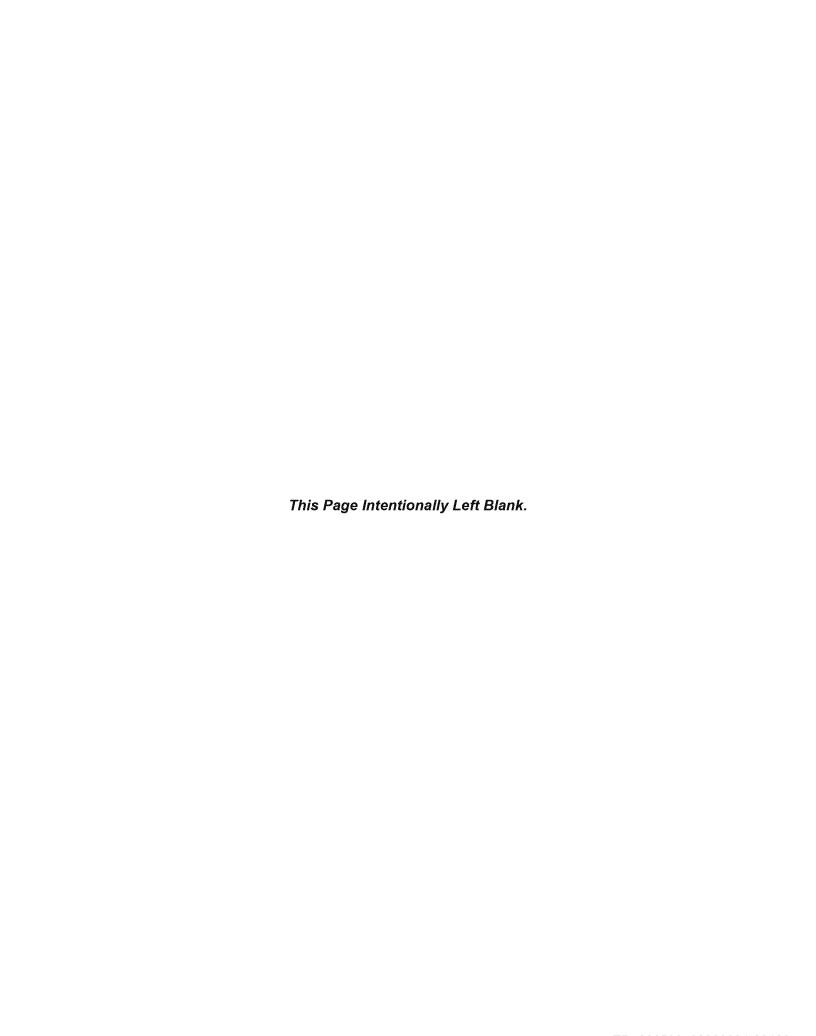


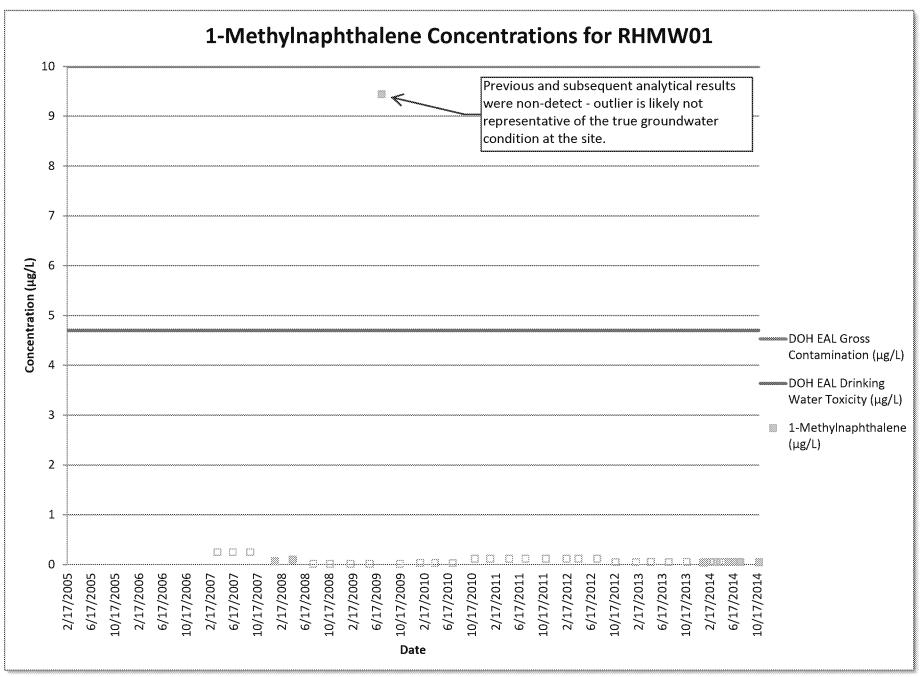


The Site-Specific Risk-Based Level (SSRBL) is 4,500 µg/L.

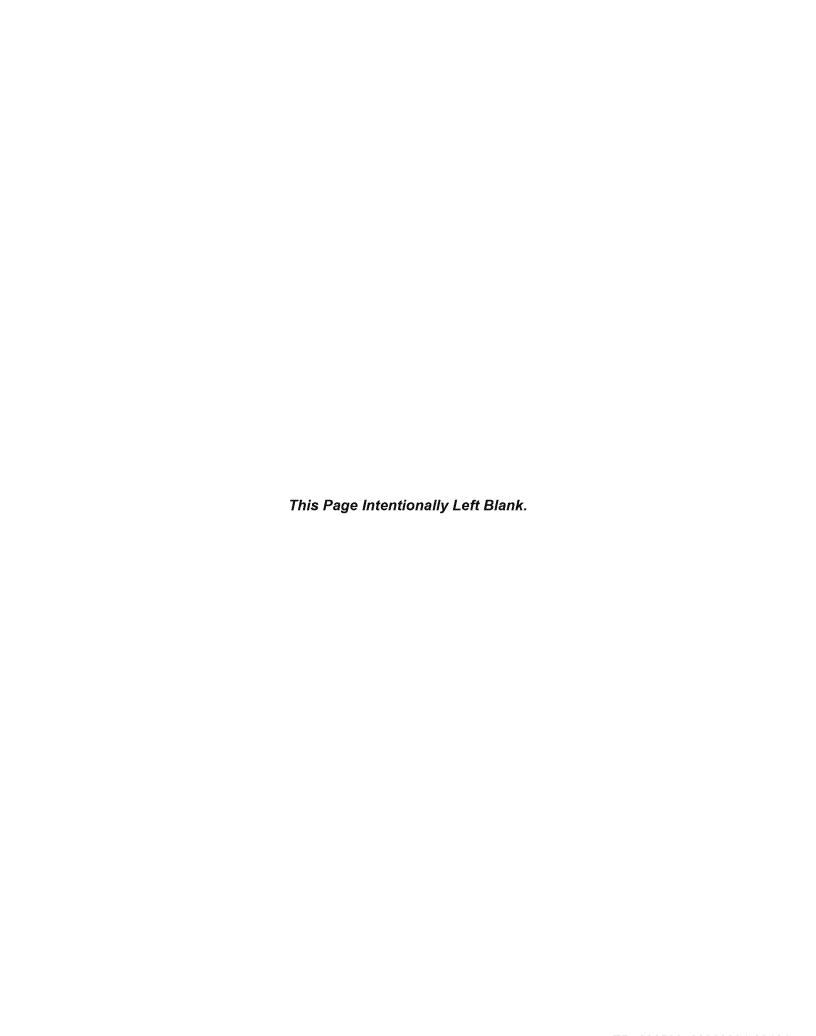
Numerous sample results had a chromatographic pattern that did not match the calibration standard. The relatively high TPH-d values may not necessarily be indicative that there is diesel fuel or other petroleum products in the well.

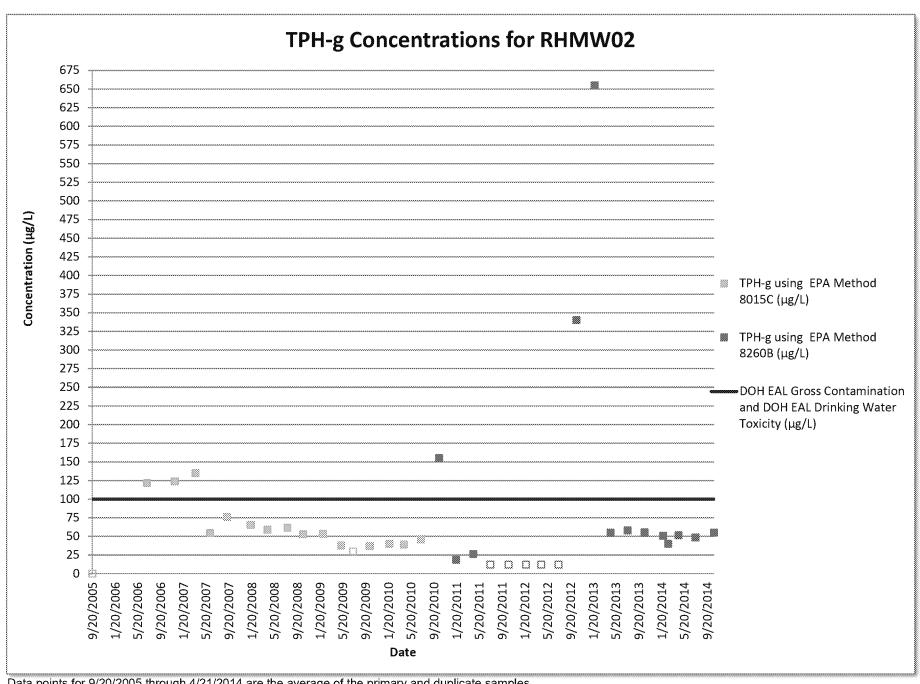
Data points for 2/17/2005 through 9/8/2005 and 12/6/2005 are the average of the primary and duplicate samples.





Unfilled boxes indicate non-detections. Method detection limits are shown.



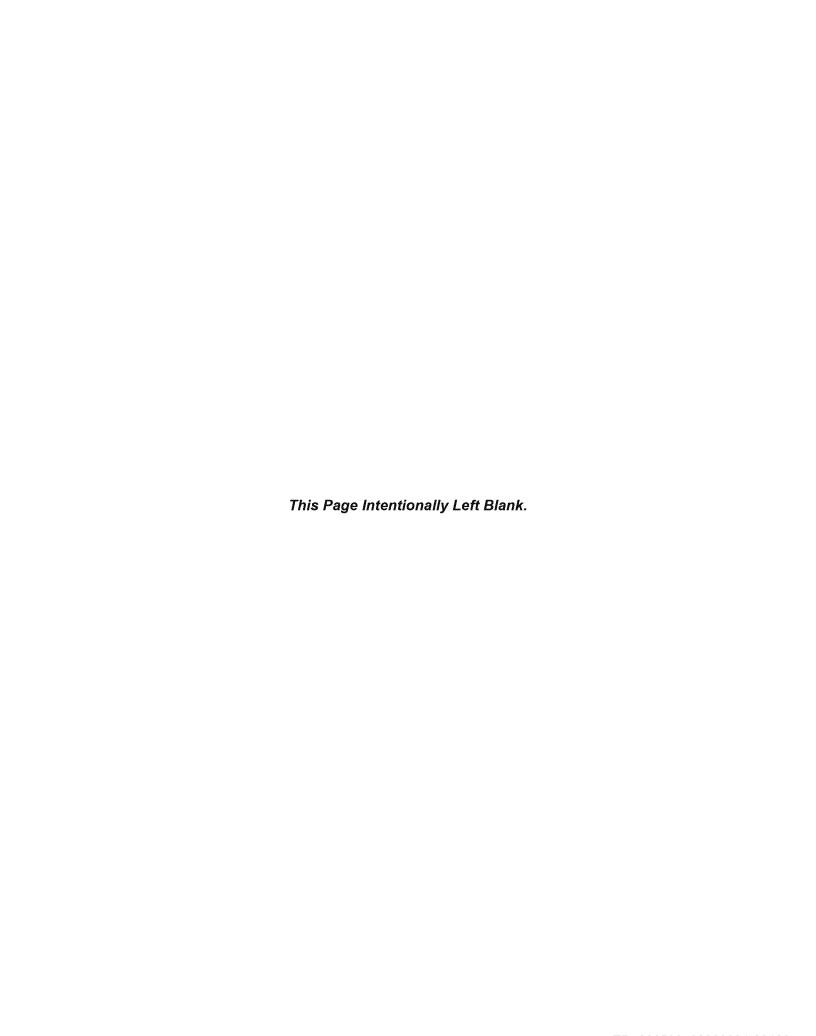


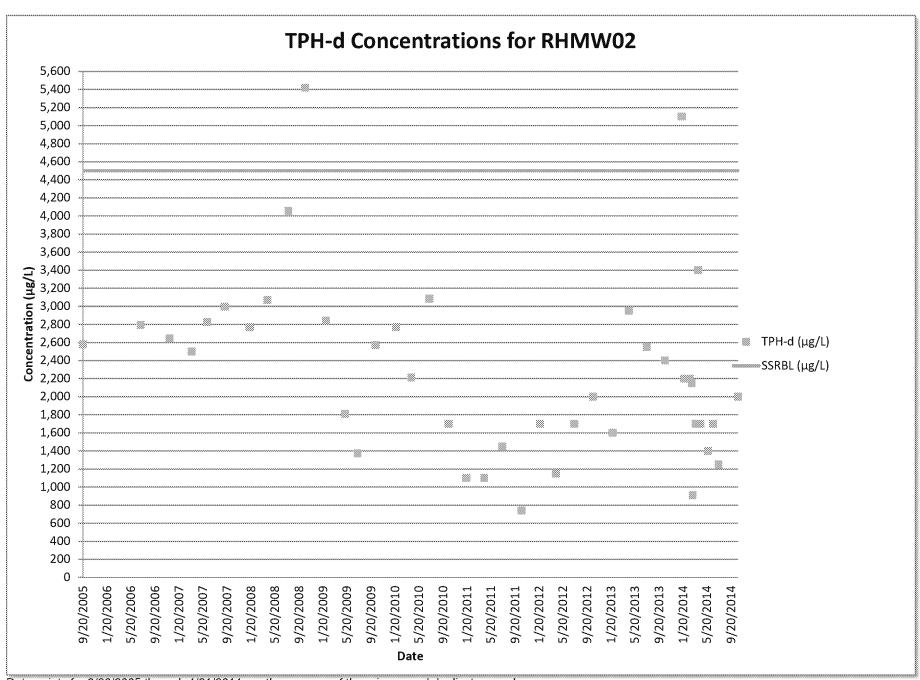
Data points for 9/20/2005 through 4/21/2014 are the average of the primary and duplicate samples.

Possible laboratory contamination for 10/21/2013 and 1/28/2014 sampling events.

Unfilled boxes indicate non-detections. Method detection limits are shown.

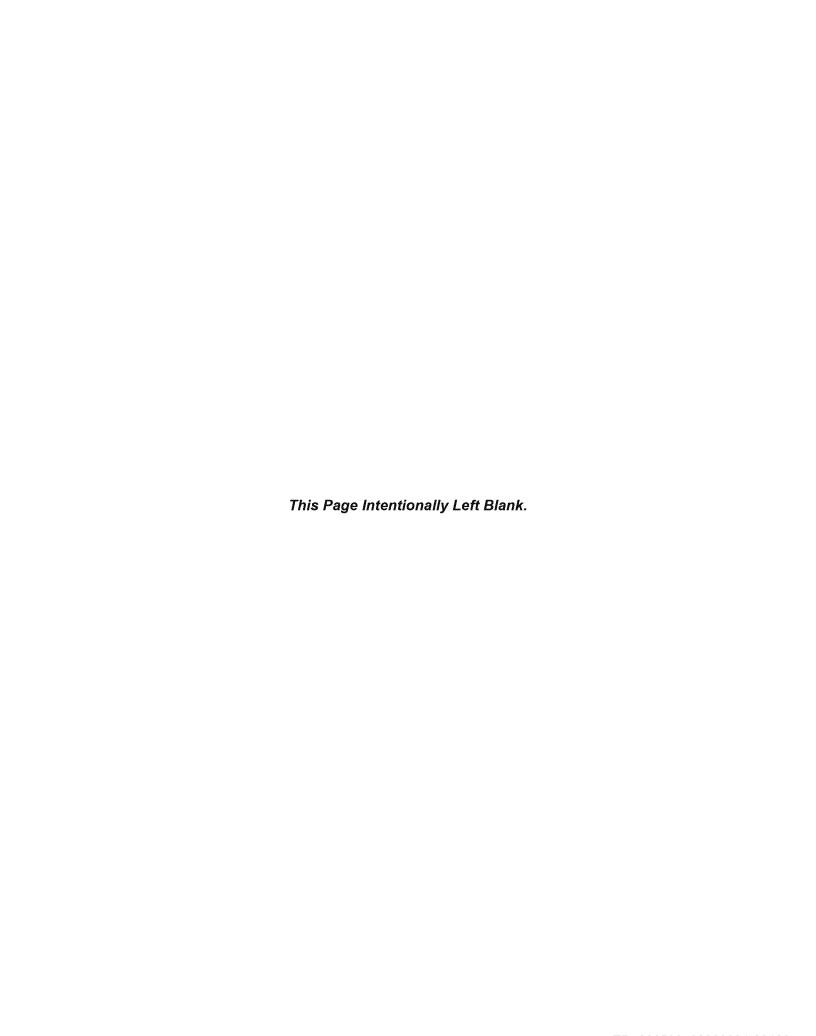
Primary sample results are shown for 1/26/2012 and 7/18/2012; all other concentrations are the average of the primary and duplicate sample results.

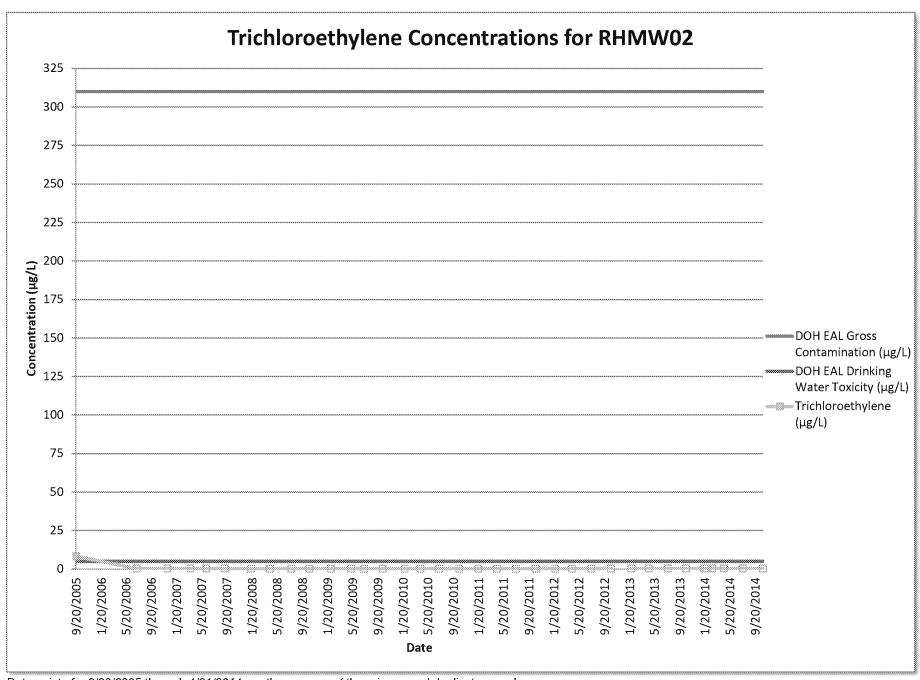




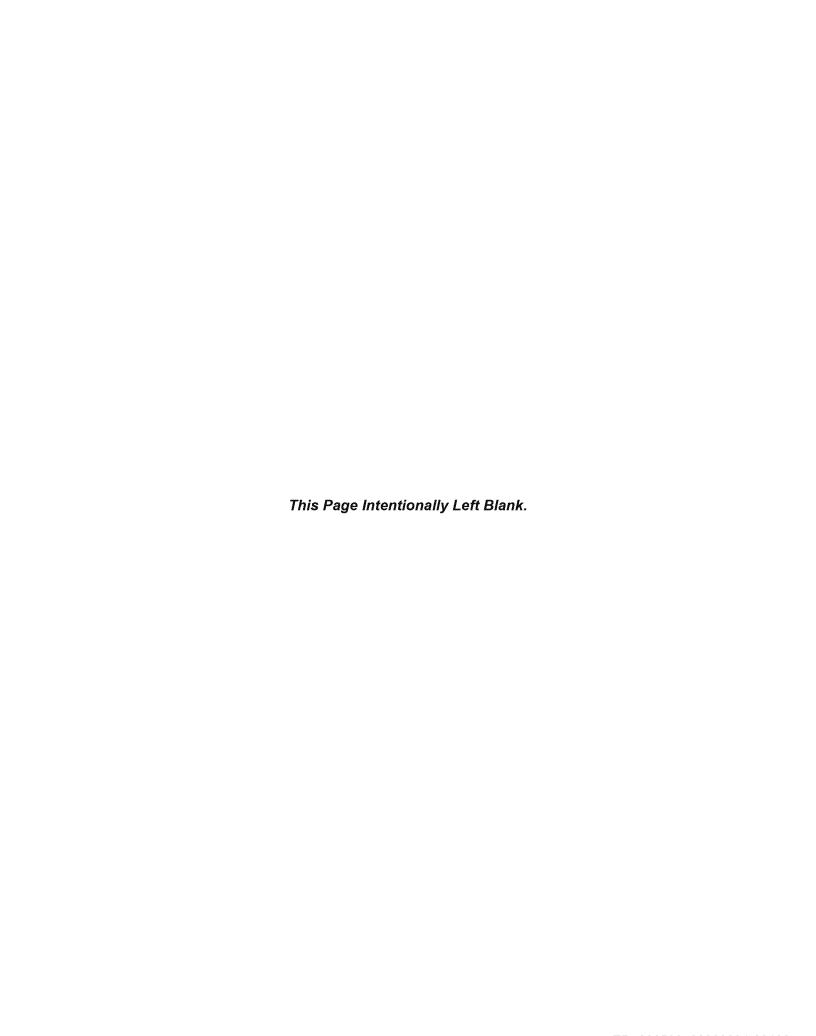
Data points for 9/20/2005 through 4/21/2014 are the average of the primary and duplicate samples.

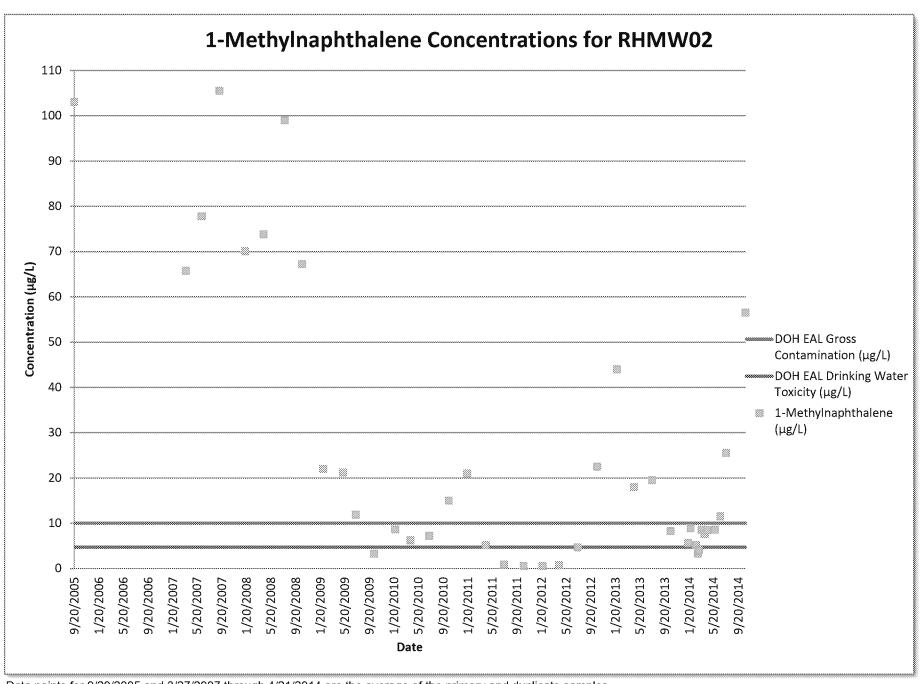
Numerous sample results had a chromatographic pattern that did not match the calibration standard. The relatively high TPH-d values may not necessarily be indicative that there is diesel fuel or other petroleum products in the well.



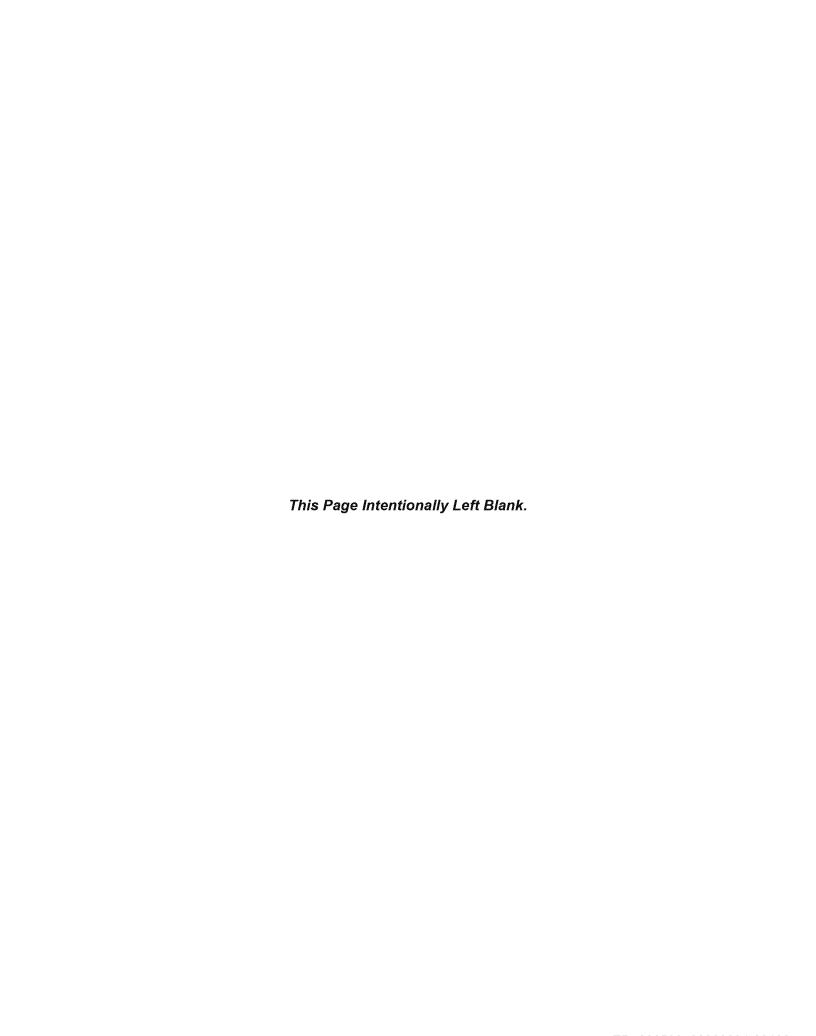


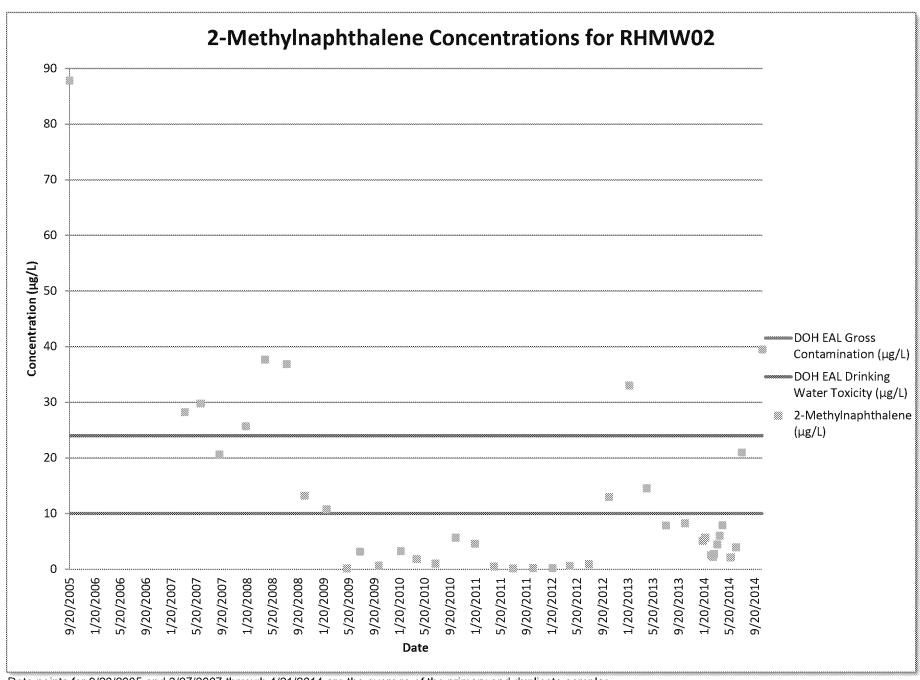
Data points for 9/20/2005 through 4/21/2014 are the average of the primary and duplicate samples. Unfilled boxes indicate non-detections. Method detection limits are shown.



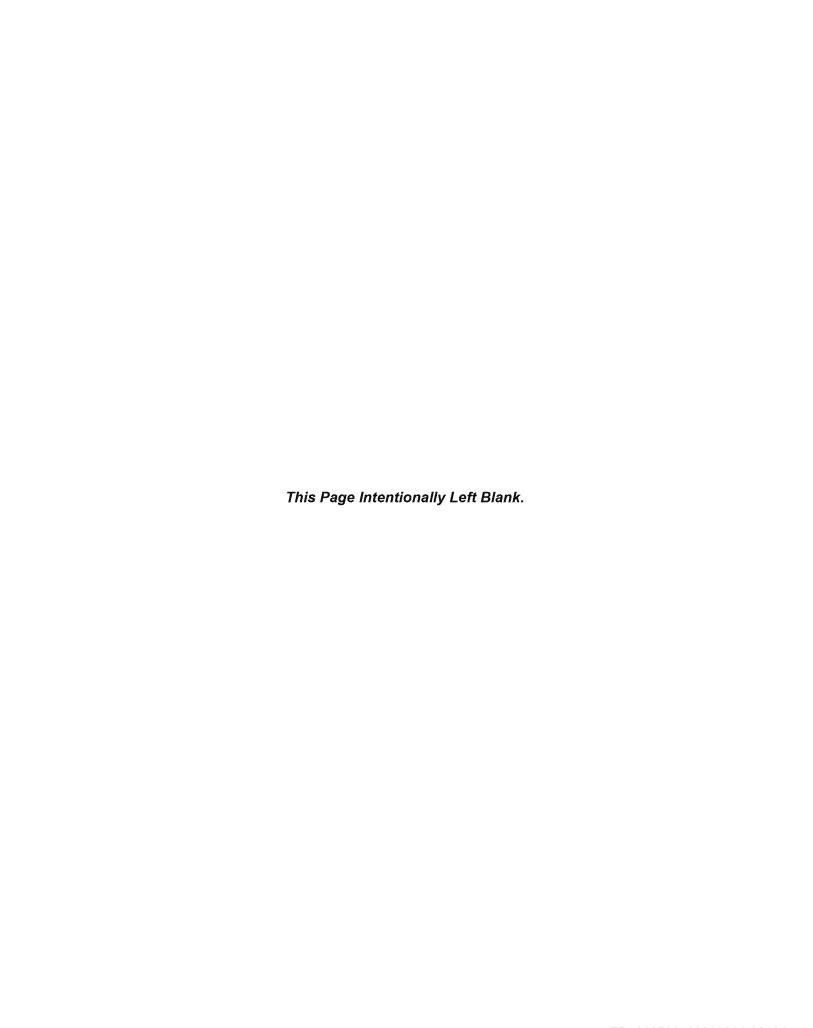


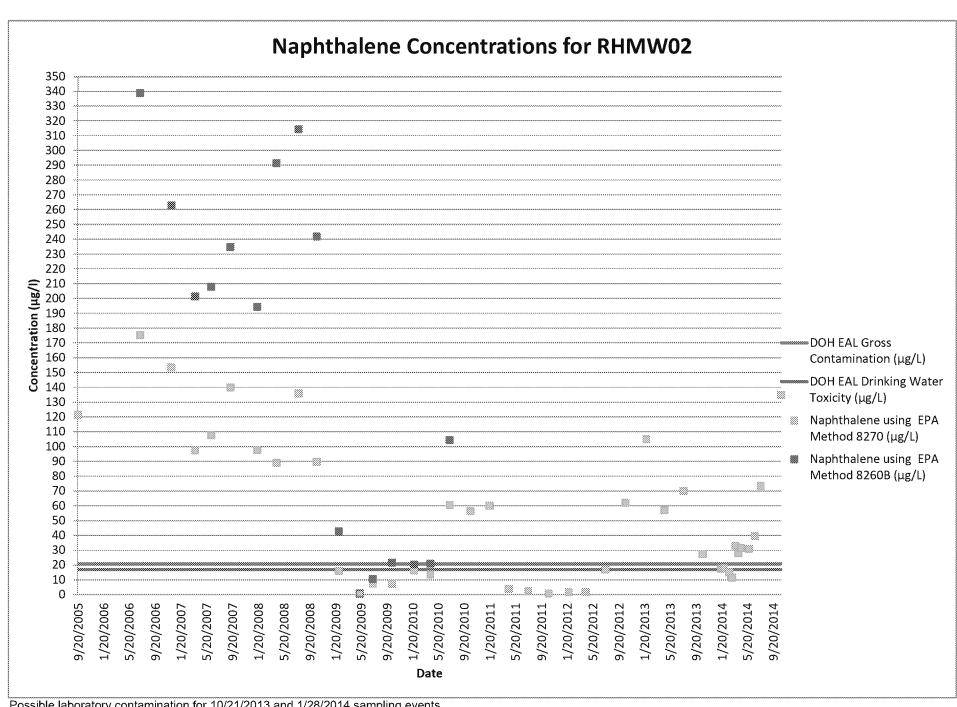
Data points for 9/20/2005 and 3/27/2007 through 4/21/2014 are the average of the primary and duplicate samples.



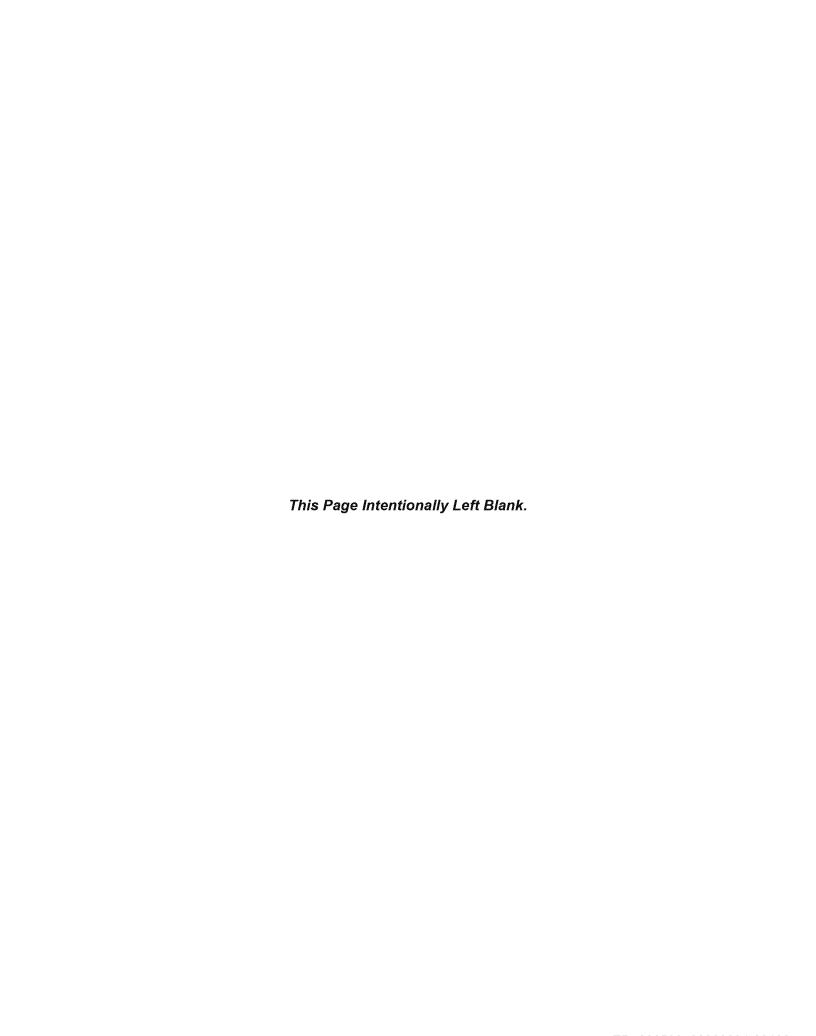


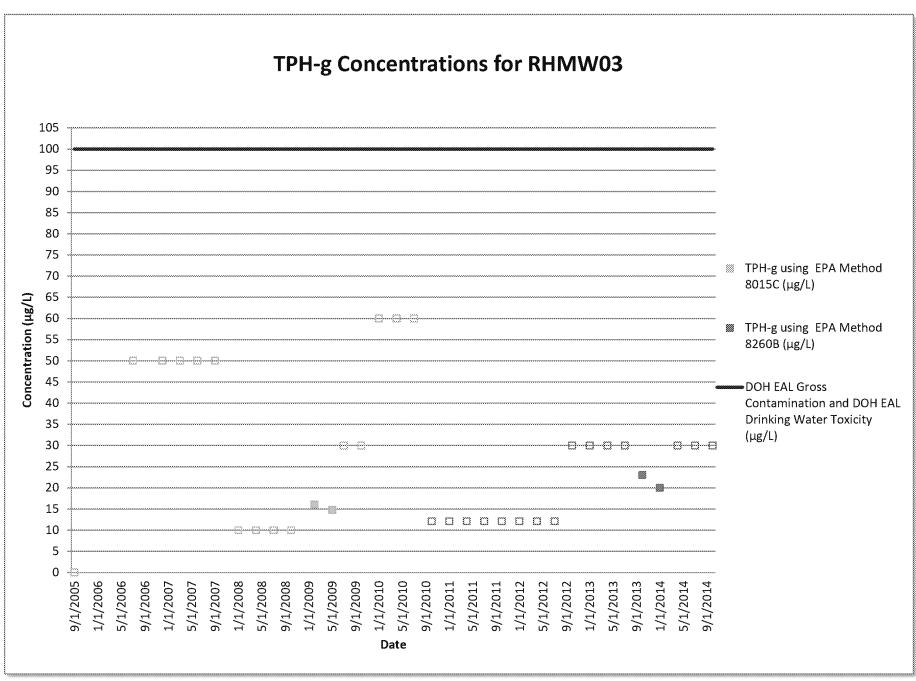
Data points for 9/20/2005 and 3/27/2007 through 4/21/2014 are the average of the primary and duplicate samples.



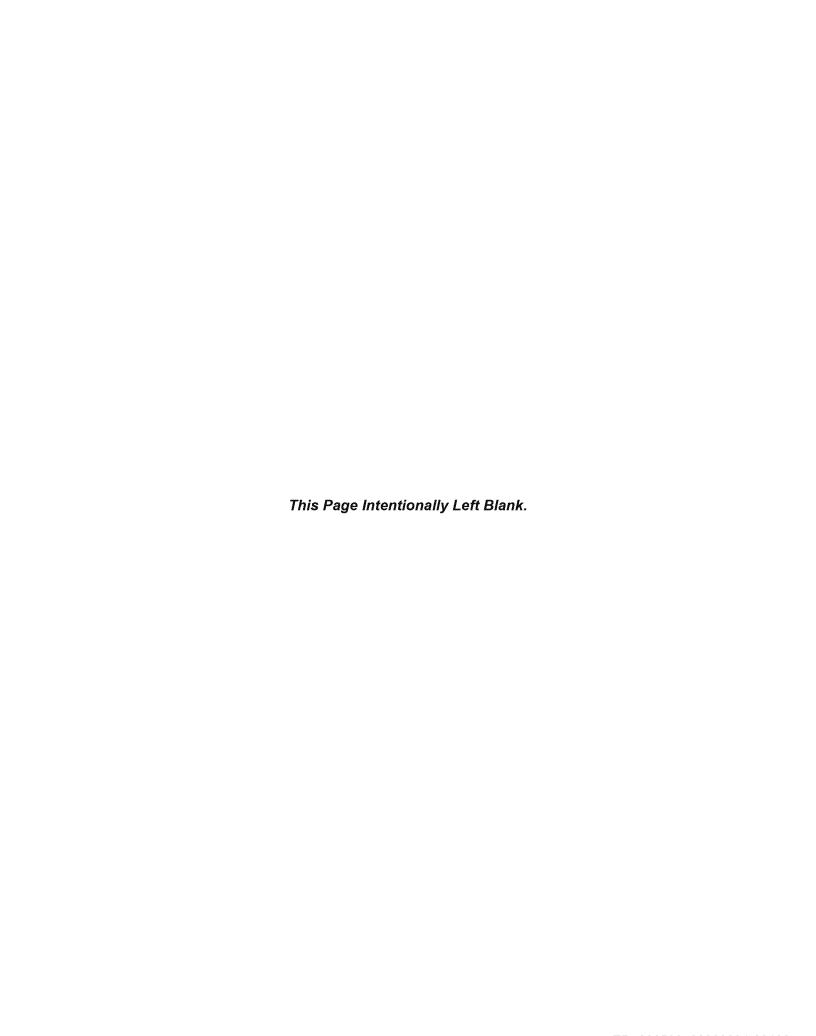


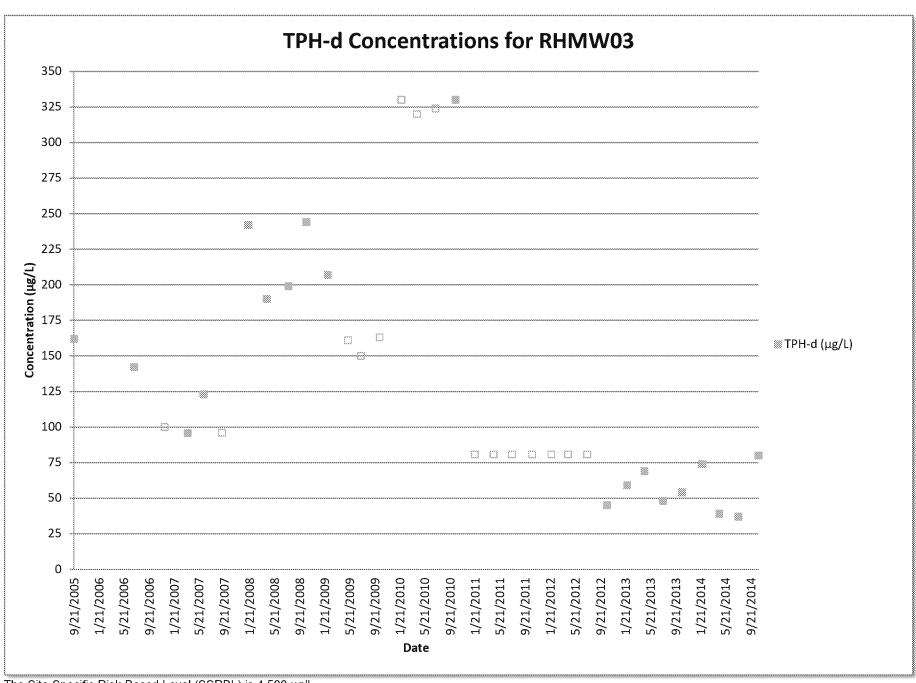
Possible laboratory contamination for 10/21/2013 and 1/28/2014 sampling events. Unfilled boxes indicate non-detections. Method detection limits are shown.





Possible laboratory contamination for 10/21/2013 and 1/28/2014 sampling events. Unfilled boxes indicate non-detections. Method detection limits are shown.

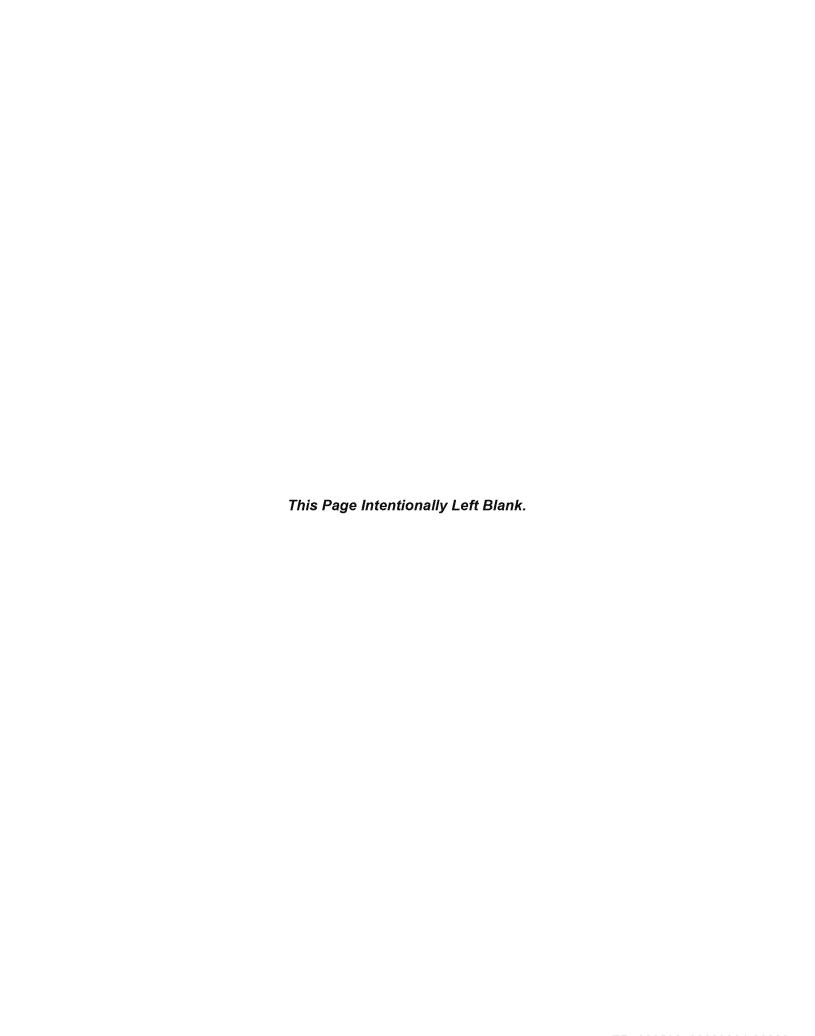


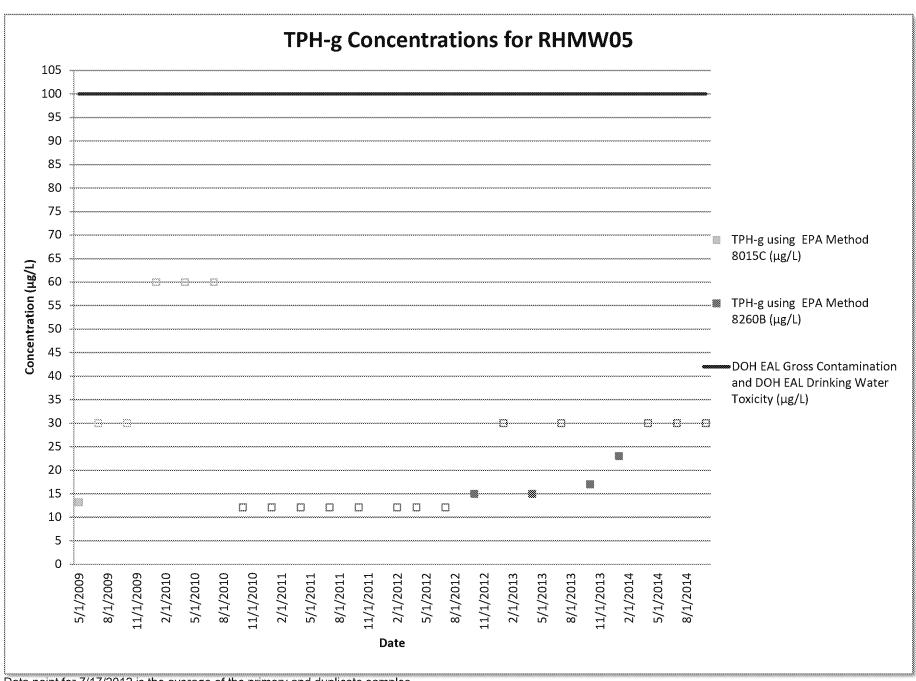


The Site-Specific Risk-Based Level (SSRBL) is 4,500 µg/L.

Unfilled boxes indicate non-detections. Method detection limits are shown.

Numerous sample results had a chromatographic pattern that didn't match the calibration standard. The relatively high TPH-d values may not necessarily be indicative that there is diesel fuel or other petroleum products in the well.

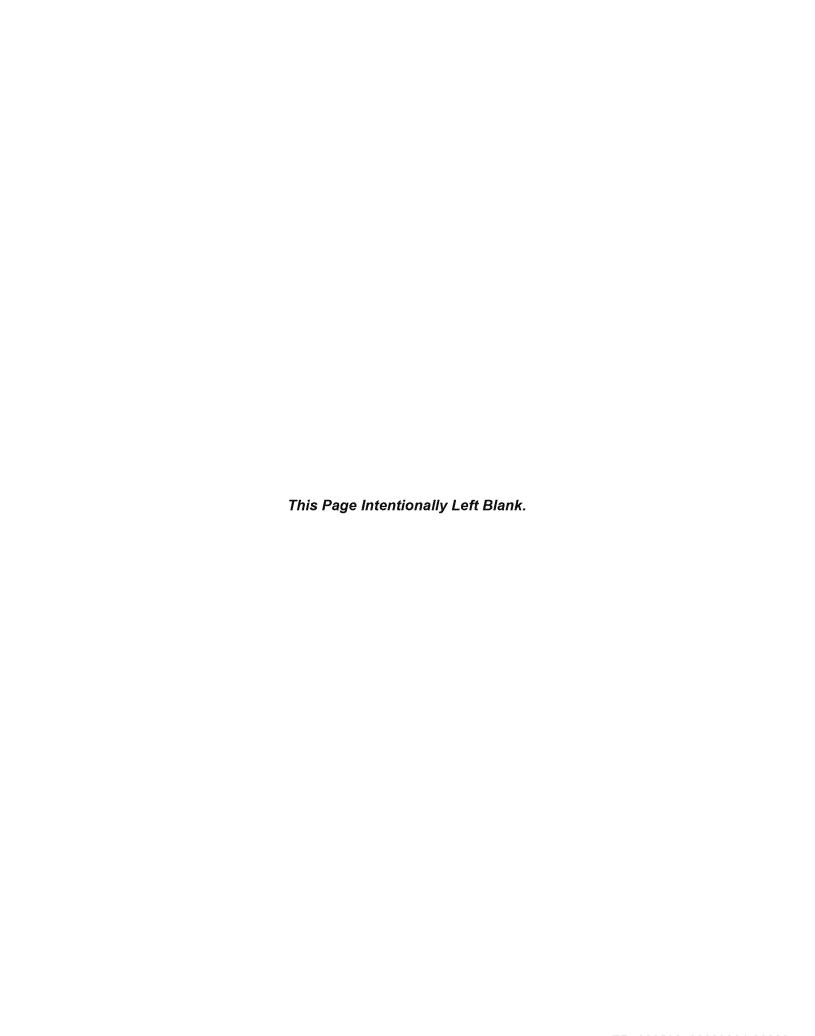


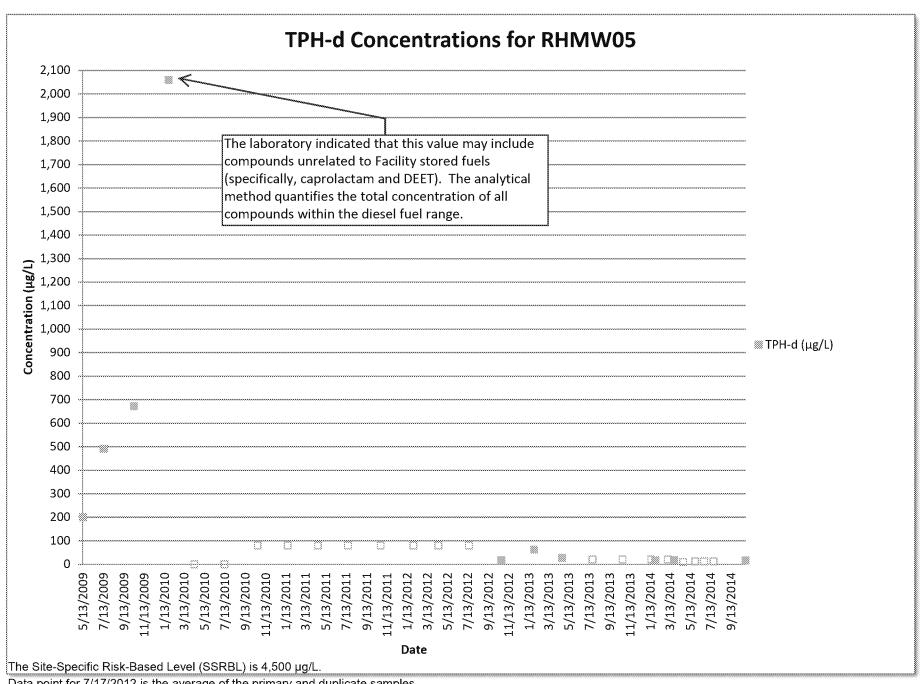


Data point for 7/17/2012 is the average of the primary and duplicate samples.

Unfilled boxes indicate non-detections. Method detection limits are shown.

Possible laboratory contamination for 10/22/2012, 10/22/2013, and 1/29/2014 sampling events.

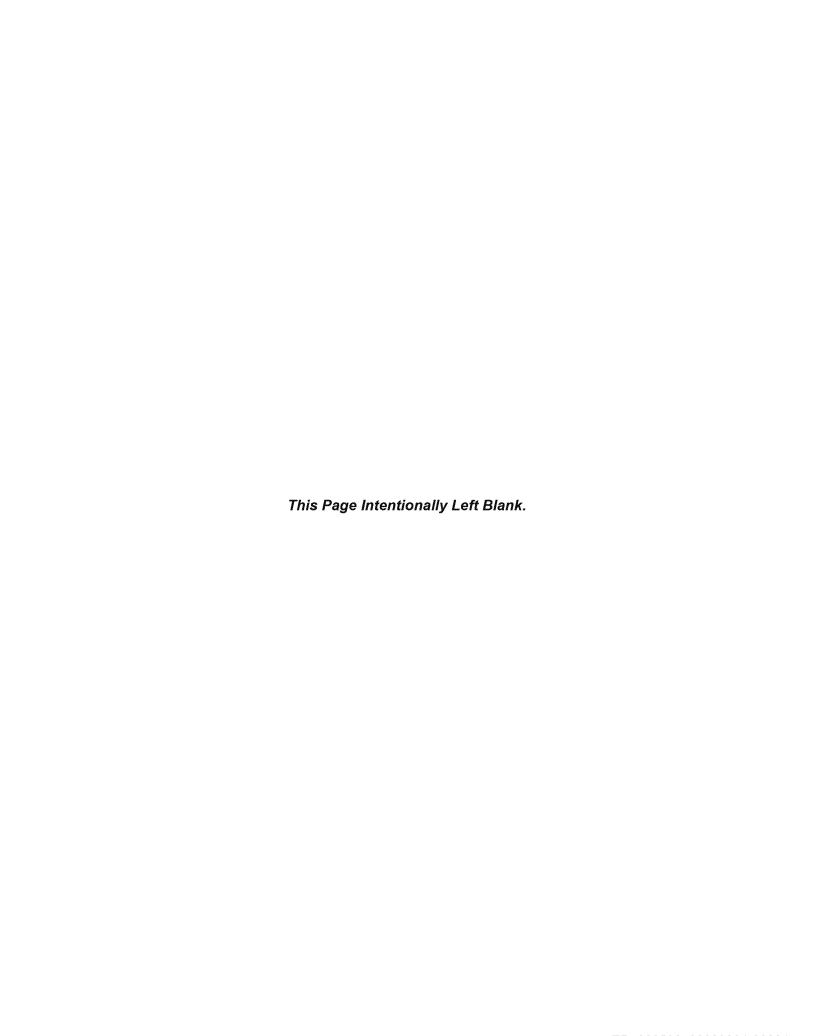


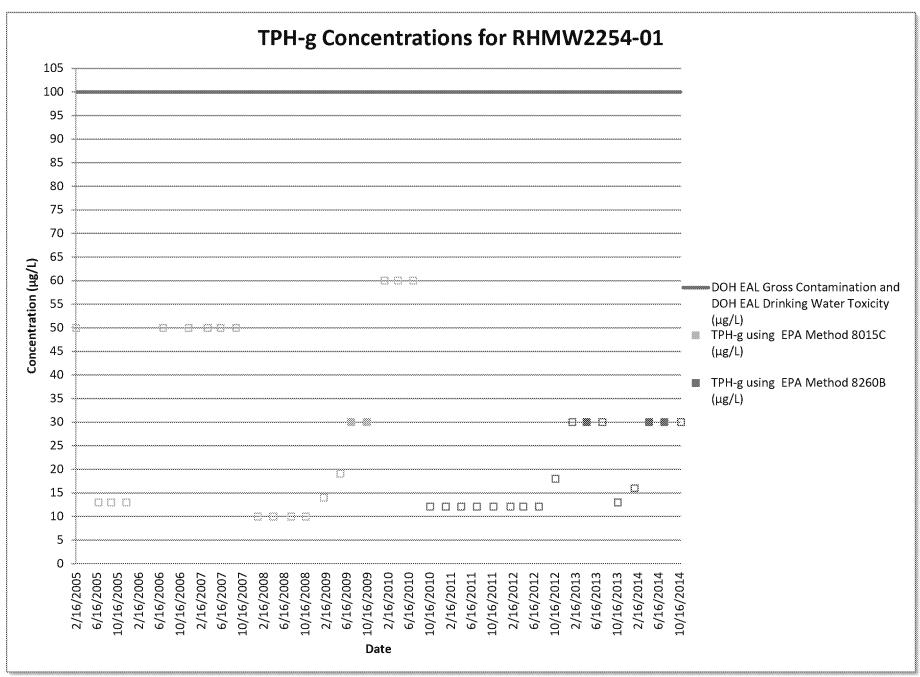


Data point for 7/17/2012 is the average of the primary and duplicate samples.

Unfilled boxes indicate non-detections. Method detection limits are shown.

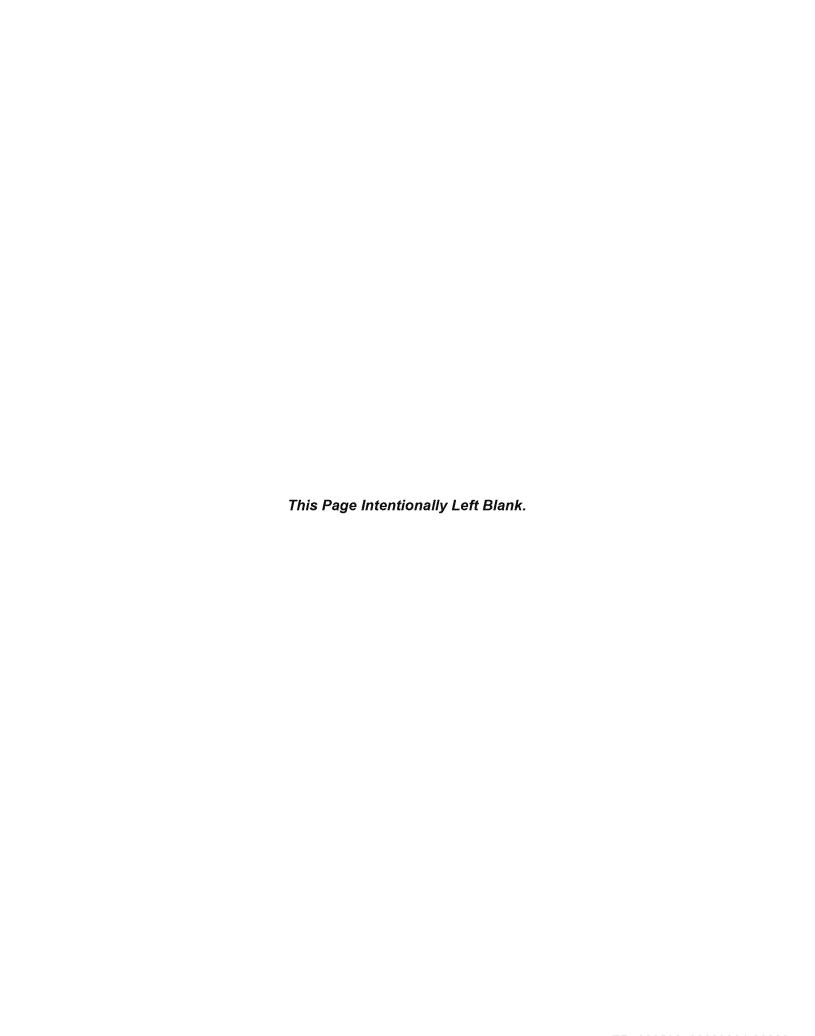
Numerous sample results had a chromatographic pattern that did not match the calibration standard. The relatively high TPH-d values may not necessarily be indicative that there is diesel fuel or other petroleum products in the well.

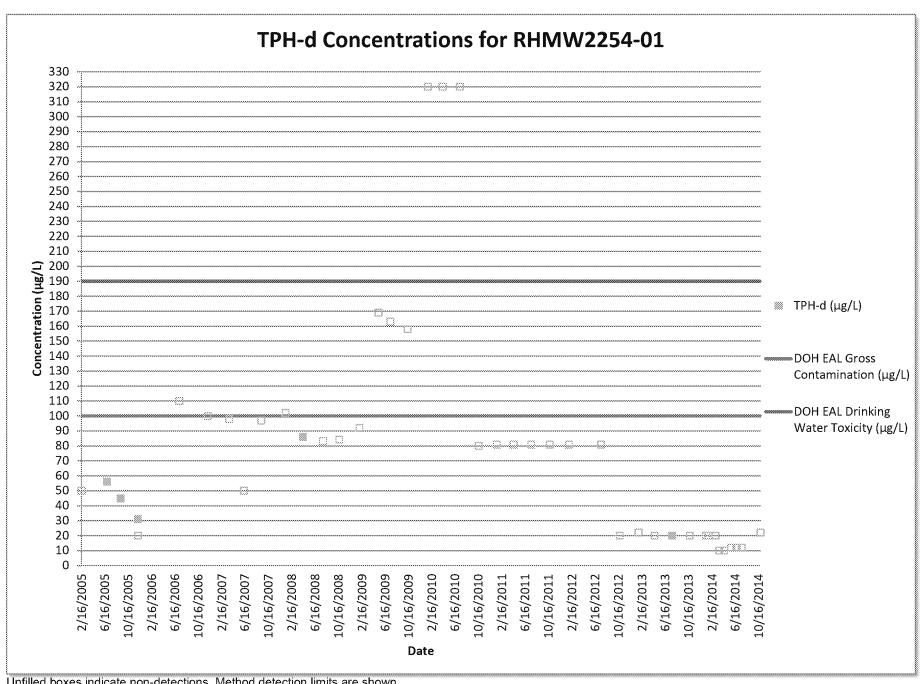




Unfilled boxes indicate non-detections. Method detection limits are shown.

Possible laboratory contamination for 10/22/2012, 10/22/2013, and 1/29/2014 sampling events.

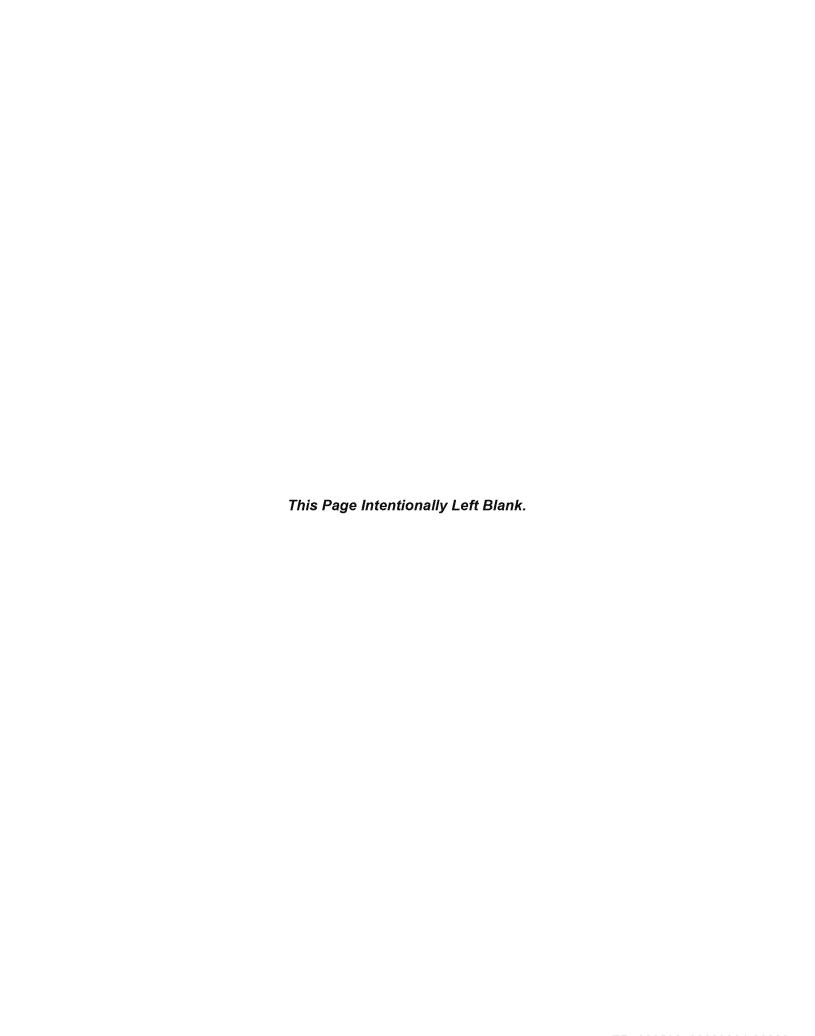




Unfilled boxes indicate non-detections. Method detection limits are shown.

Laboratory data rejected for 1/15/2008 sampling event.

Numerous sample results had a chromatographic pattern that did not match the calibration standard. The relatively high TPH-d values may not necessarily be indicative that there is diesel fuel or other petroleum products in the well.



APPENDIX E

Cumulative Groundwater Results (included on attached CD)

